



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 08:57 am BST

PDB ID : 5XFM
Title : Crystal structure of beta-arabinopyranosidase
Authors : Kato, K.; Okuyama, M.; Yao, M.
Deposited on : 2017-04-10
Resolution : 2.30 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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

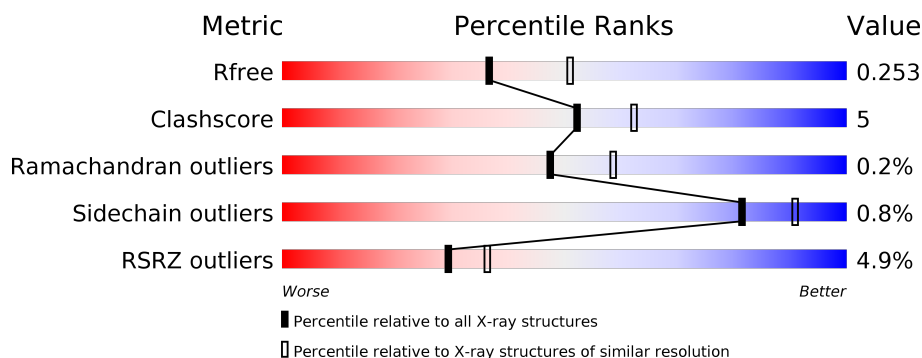
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	642	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>14%</div> <div>•</div> </div> </div>
1	B	642	<div> <div>5%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>•</div> </div> </div>
1	C	642	<div> <div>5%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>•</div> </div> </div>
1	D	642	<div> <div>8%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>••</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21423 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 Alpha-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	632	Total	C	N	O	S	0	0	0
			5056	3204	870	953	29			
1	B	628	Total	C	N	O	S	0	0	0
			5021	3182	864	947	28			
1	C	632	Total	C	N	O	S	0	0	0
			5056	3204	870	953	29			
1	D	629	Total	C	N	O	S	0	0	0
			5029	3188	865	948	28			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	MET	-	expression tag	UNP A0A1H5YQC4
A	7	ASN	-	expression tag	UNP A0A1H5YQC4
A	8	HIS	-	expression tag	UNP A0A1H5YQC4
A	9	LYS	-	expression tag	UNP A0A1H5YQC4
A	10	VAL	-	expression tag	UNP A0A1H5YQC4
A	11	HIS	-	expression tag	UNP A0A1H5YQC4
A	12	HIS	-	expression tag	UNP A0A1H5YQC4
A	13	HIS	-	expression tag	UNP A0A1H5YQC4
A	14	HIS	-	expression tag	UNP A0A1H5YQC4
A	15	HIS	-	expression tag	UNP A0A1H5YQC4
A	16	HIS	-	expression tag	UNP A0A1H5YQC4
A	17	MET	-	expression tag	UNP A0A1H5YQC4
A	18	GLU	-	expression tag	UNP A0A1H5YQC4
A	19	LEU	-	expression tag	UNP A0A1H5YQC4
A	498	LEU	PHE	engineered mutation	UNP A0A1H5YQC4
B	6	MET	-	expression tag	UNP A0A1H5YQC4
B	7	ASN	-	expression tag	UNP A0A1H5YQC4
B	8	HIS	-	expression tag	UNP A0A1H5YQC4
B	9	LYS	-	expression tag	UNP A0A1H5YQC4
B	10	VAL	-	expression tag	UNP A0A1H5YQC4
B	11	HIS	-	expression tag	UNP A0A1H5YQC4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	12	HIS	-	expression tag	UNP A0A1H5YQC4
B	13	HIS	-	expression tag	UNP A0A1H5YQC4
B	14	HIS	-	expression tag	UNP A0A1H5YQC4
B	15	HIS	-	expression tag	UNP A0A1H5YQC4
B	16	HIS	-	expression tag	UNP A0A1H5YQC4
B	17	MET	-	expression tag	UNP A0A1H5YQC4
B	18	GLU	-	expression tag	UNP A0A1H5YQC4
B	19	LEU	-	expression tag	UNP A0A1H5YQC4
B	498	LEU	PHE	engineered mutation	UNP A0A1H5YQC4
C	6	MET	-	expression tag	UNP A0A1H5YQC4
C	7	ASN	-	expression tag	UNP A0A1H5YQC4
C	8	HIS	-	expression tag	UNP A0A1H5YQC4
C	9	LYS	-	expression tag	UNP A0A1H5YQC4
C	10	VAL	-	expression tag	UNP A0A1H5YQC4
C	11	HIS	-	expression tag	UNP A0A1H5YQC4
C	12	HIS	-	expression tag	UNP A0A1H5YQC4
C	13	HIS	-	expression tag	UNP A0A1H5YQC4
C	14	HIS	-	expression tag	UNP A0A1H5YQC4
C	15	HIS	-	expression tag	UNP A0A1H5YQC4
C	16	HIS	-	expression tag	UNP A0A1H5YQC4
C	17	MET	-	expression tag	UNP A0A1H5YQC4
C	18	GLU	-	expression tag	UNP A0A1H5YQC4
C	19	LEU	-	expression tag	UNP A0A1H5YQC4
C	498	LEU	PHE	engineered mutation	UNP A0A1H5YQC4
D	6	MET	-	expression tag	UNP A0A1H5YQC4
D	7	ASN	-	expression tag	UNP A0A1H5YQC4
D	8	HIS	-	expression tag	UNP A0A1H5YQC4
D	9	LYS	-	expression tag	UNP A0A1H5YQC4
D	10	VAL	-	expression tag	UNP A0A1H5YQC4
D	11	HIS	-	expression tag	UNP A0A1H5YQC4
D	12	HIS	-	expression tag	UNP A0A1H5YQC4
D	13	HIS	-	expression tag	UNP A0A1H5YQC4
D	14	HIS	-	expression tag	UNP A0A1H5YQC4
D	15	HIS	-	expression tag	UNP A0A1H5YQC4
D	16	HIS	-	expression tag	UNP A0A1H5YQC4
D	17	MET	-	expression tag	UNP A0A1H5YQC4
D	18	GLU	-	expression tag	UNP A0A1H5YQC4
D	19	LEU	-	expression tag	UNP A0A1H5YQC4
D	498	LEU	PHE	engineered mutation	UNP A0A1H5YQC4

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

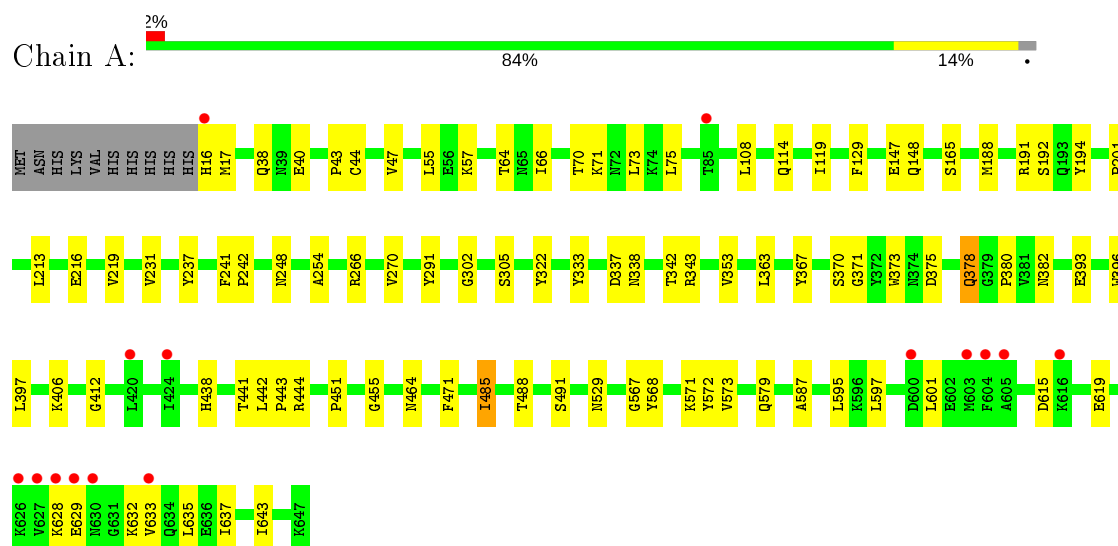
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	333	Total O 333 333	0	0
3	B	291	Total O 291 291	0	0
3	C	349	Total O 349 349	0	0
3	D	284	Total O 284 284	0	0

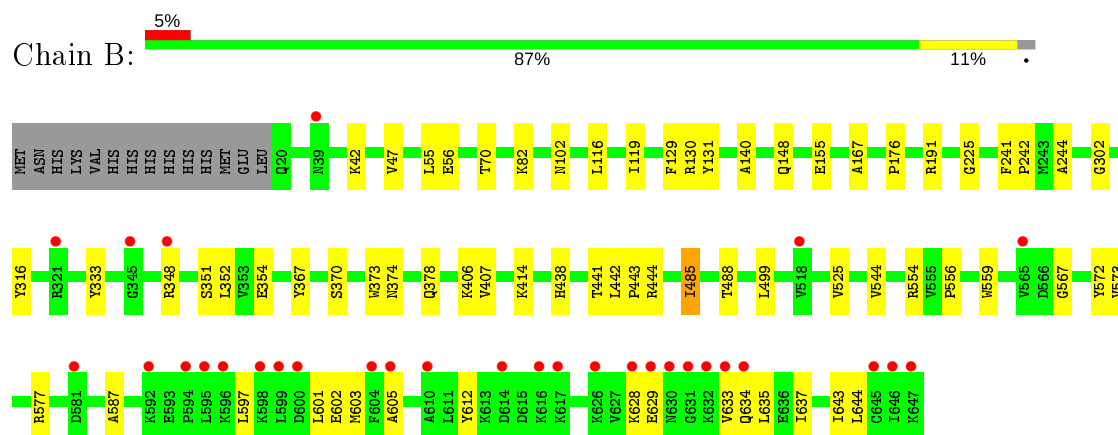
3 Residue-property plots [i](#)

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.

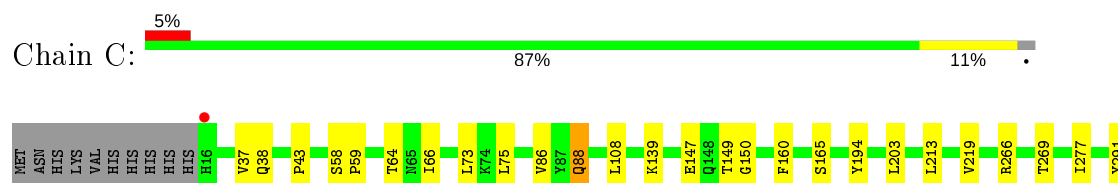
• Molecule 1: Alpha-glucosidase

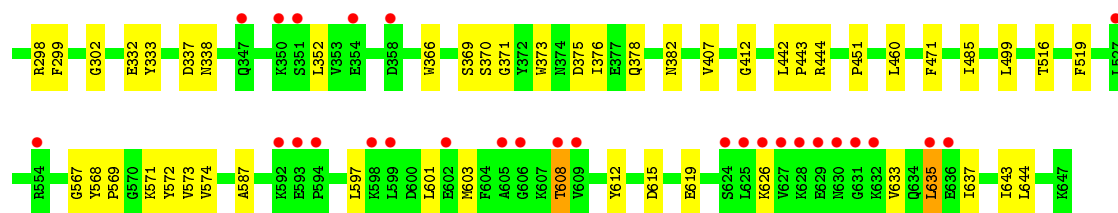


• Molecule 1: Alpha-glucosidase

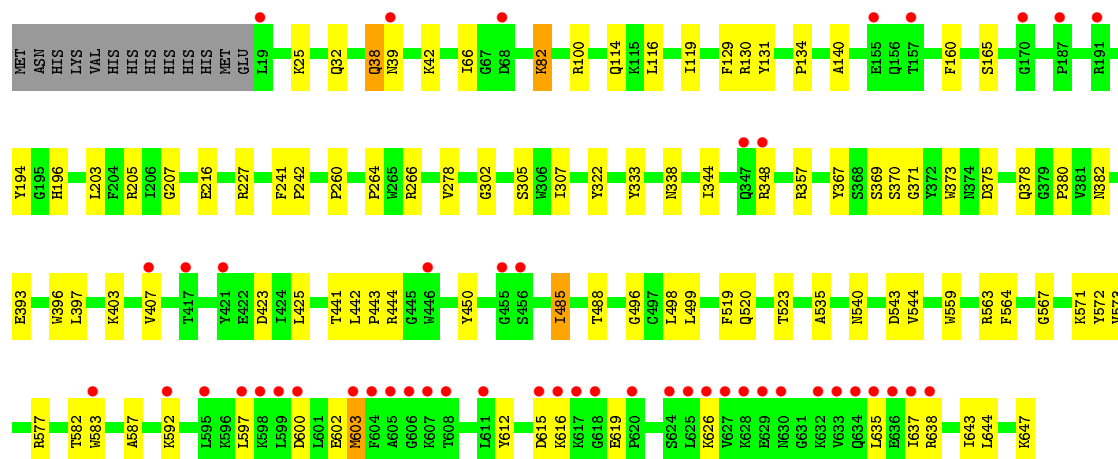
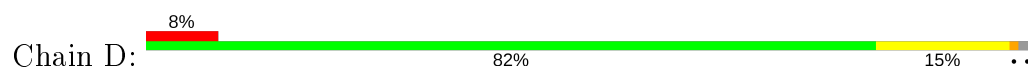


• Molecule 1: Alpha-glucosidase





● Molecule 1: Alpha-glucosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	72.81Å 90.51Å 128.55Å 105.50° 94.80° 96.15°	Depositor
Resolution (Å)	45.78 – 2.30 45.78 – 2.30	Depositor EDS
% Data completeness (in resolution range)	90.0 (45.78-2.30) 90.1 (45.78-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.207 , 0.253 0.209 , 0.253	Depositor DCC
R_{free} test set	6269 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.567	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.6	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	0.92	EDS
Total number of atoms	21423	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 5.77% 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: CA

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	0.31	0/5180	0.53	0/7022
1	B	0.30	0/5144	0.51	0/6974
1	C	0.32	0/5180	0.53	1/7022 (0.0%)
1	D	0.30	0/5152	0.52	0/6985
All	All	0.31	0/20656	0.52	1/28003 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	635	LEU	CA-CB-CG	6.38	129.98	115.30

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	5056	0	4912	62	0
1	B	5021	0	4879	42	0
1	C	5056	0	4912	51	0
1	D	5029	0	4890	65	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	333	0	0	7	0
3	B	291	0	0	7	0
3	C	349	0	0	6	0
3	D	284	0	0	9	0
All	All	21423	0	19593	217	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 (217) 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:64:THR:HG22	1:C:66:ILE:H	1.44	0.82
1:A:70:THR:HG22	1:A:71:LYS:HG3	1.62	0.81
1:D:496:GLY:O	3:D:801:HOH:O	1.98	0.80
1:D:583:TRP:O	3:D:802:HOH:O	2.00	0.79
1:D:344:ILE:HA	1:D:348:ARG:HH21	1.48	0.77
1:A:191:ARG:NH1	1:A:192:SER:O	2.16	0.77
1:D:160:PHE:HB2	1:D:203:LEU:HB3	1.68	0.75
1:A:64:THR:HG22	1:A:66:ILE:H	1.52	0.72
1:A:38:GLN:OE1	1:A:75:LEU:N	2.21	0.71
1:D:615:ASP:O	1:D:638:ARG:NH2	2.25	0.70
1:D:338:ASN:HD22	1:D:378:GLN:HE21	1.38	0.69
1:C:38:GLN:OE1	1:C:75:LEU:N	2.23	0.67
1:C:635:LEU:HD23	1:C:643:ILE:HD13	1.80	0.64
1:B:637:ILE:HG12	1:B:643:ILE:HD12	1.80	0.64
1:C:43:PRO:HG3	1:C:73:LEU:HD22	1.79	0.63
1:D:647:LYS:N	3:D:802:HOH:O	2.31	0.63
1:A:637:ILE:HG12	1:A:643:ILE:HD12	1.81	0.63
1:A:338:ASN:ND2	1:A:378:GLN:OE1	2.32	0.63
1:A:441:THR:HB	3:A:803:HOH:O	1.99	0.63
1:A:635:LEU:HD13	1:A:643:ILE:HG21	1.81	0.62
1:B:370:SER:HA	1:B:373:TRP:CZ2	2.34	0.62
1:A:396:TRP:CE3	1:A:397:LEU:HD12	2.35	0.61
1:A:628:LYS:HD2	1:A:629:GLU:H	1.65	0.61
1:A:147:GLU:OE2	1:A:266:ARG:NH2	2.32	0.60
1:A:43:PRO:HG3	1:A:73:LEU:HD22	1.82	0.60
1:D:396:TRP:CE3	1:D:397:LEU:HD12	2.37	0.60
1:D:25:LYS:HG2	1:D:32:GLN:HG2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:499:LEU:O	3:C:801:HOH:O	2.16	0.59
1:A:254:ALA:O	3:A:802:HOH:O	2.16	0.59
1:B:56:GLU:OE2	3:B:802:HOH:O	2.17	0.59
1:D:423:ASP:OD1	3:D:803:HOH:O	2.17	0.59
1:C:376:ILE:HG23	1:C:378:GLN:HG2	1.85	0.58
1:A:579:GLN:NE2	3:A:811:HOH:O	2.36	0.58
1:C:637:ILE:HG12	1:C:643:ILE:HD12	1.85	0.58
1:A:73:LEU:HD23	1:A:108:LEU:HD13	1.85	0.57
1:C:370:SER:HA	1:C:373:TRP:CZ2	2.39	0.57
1:A:396:TRP:HE3	1:A:397:LEU:HD12	1.70	0.57
1:A:370:SER:HA	1:A:373:TRP:CZ2	2.39	0.57
1:A:231:VAL:HG22	1:A:237:TYR:CZ	2.39	0.56
1:B:70:THR:HG23	1:B:148:GLN:OE1	2.05	0.56
1:B:244:ALA:O	1:B:414:LYS:NZ	2.38	0.56
1:A:491:SER:HB3	1:A:529:ASN:ND2	2.20	0.56
1:B:499:LEU:HD13	1:B:544:VAL:HG11	1.89	0.55
1:D:393:GLU:O	1:D:397:LEU:HD13	2.06	0.55
1:A:70:THR:HG23	1:A:148:GLN:OE1	2.07	0.54
1:C:635:LEU:HD23	1:C:643:ILE:HG21	1.90	0.54
1:C:442:LEU:HD12	1:C:443:PRO:HD2	1.89	0.54
1:B:348:ARG:NH1	3:B:827:HOH:O	2.41	0.54
1:A:231:VAL:HG22	1:A:237:TYR:CE2	2.42	0.54
1:C:165:SER:HA	1:C:194:TYR:CD1	2.43	0.53
1:B:378:GLN:NE2	3:B:824:HOH:O	2.40	0.53
1:C:147:GLU:OE2	1:C:266:ARG:NH2	2.38	0.53
1:D:196:HIS:O	1:D:227:ARG:NH1	2.41	0.53
1:B:155:GLU:OE1	3:B:802:HOH:O	2.19	0.53
1:A:191:ARG:HH11	1:A:191:ARG:HG3	1.72	0.53
1:D:370:SER:HA	1:D:373:TRP:CZ2	2.43	0.53
1:D:540:ASN:HA	1:D:543:ASP:OD1	2.09	0.53
1:D:442:LEU:HD12	1:D:443:PRO:HD2	1.91	0.52
1:C:338:ASN:ND2	1:C:378:GLN:HE21	2.08	0.52
1:A:375:ASP:HA	1:B:140:ALA:HB3	1.91	0.52
1:A:412:GLY:O	3:A:803:HOH:O	2.18	0.51
1:C:149:THR:HG21	1:C:266:ARG:HD2	1.91	0.51
1:A:353:VAL:HG22	1:A:363:LEU:HD12	1.91	0.51
1:B:577:ARG:NH2	3:B:831:HOH:O	2.44	0.51
1:C:633:VAL:HG12	1:C:635:LEU:CD1	2.41	0.51
1:C:160:PHE:HB2	1:C:203:LEU:HB3	1.92	0.51
1:A:188:MET:O	1:A:231:VAL:HG21	2.11	0.50
1:D:116:LEU:HD11	1:D:131:TYR:HB3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:563:ARG:NH1	1:D:602:GLU:OE2	2.44	0.50
1:D:637:ILE:HG12	1:D:643:ILE:HD12	1.93	0.50
1:B:525:VAL:O	1:B:577:ARG:NH1	2.45	0.50
1:B:47:VAL:HB	1:B:55:LEU:HB2	1.92	0.50
1:D:571:LYS:NZ	3:D:817:HOH:O	2.37	0.50
1:D:635:LEU:HD13	1:D:643:ILE:HG21	1.94	0.50
1:B:348:ARG:HH11	1:B:348:ARG:HG2	1.77	0.49
1:A:165:SER:HA	1:A:194:TYR:CD1	2.47	0.49
1:B:601:LEU:HD12	1:B:633:VAL:HG21	1.93	0.49
1:A:406:LYS:HE3	1:A:438:HIS:CD2	2.48	0.49
1:C:516:THR:O	1:C:519:PHE:HB3	2.13	0.49
1:A:219:VAL:HB	1:A:444:ARG:HD3	1.95	0.49
1:C:635:LEU:CD2	1:C:643:ILE:HG21	2.43	0.49
1:A:393:GLU:O	1:A:397:LEU:HD13	2.12	0.49
1:D:603:MET:HG3	1:D:603:MET:H	1.39	0.48
1:A:601:LEU:HD12	1:A:633:VAL:HG21	1.96	0.48
1:D:378:GLN:HG3	3:D:830:HOH:O	2.14	0.48
1:B:82:LYS:HG2	1:B:102:ASN:OD1	2.14	0.48
1:C:302:GLY:HA2	1:C:333:TYR:O	2.14	0.48
1:D:38:GLN:OE1	1:D:38:GLN:HA	2.14	0.48
1:B:602:GLU:CD	1:B:602:GLU:H	2.17	0.47
1:B:634:GLN:HG2	1:B:635:LEU:N	2.29	0.47
1:A:57:LYS:NZ	3:A:801:HOH:O	1.97	0.47
1:C:298:ARG:HG2	3:C:802:HOH:O	2.13	0.47
1:A:47:VAL:HB	1:A:55:LEU:HB2	1.97	0.47
1:A:371:GLY:HA2	1:A:382:ASN:OD1	2.14	0.47
1:C:338:ASN:HD22	1:C:378:GLN:HE21	1.63	0.47
1:D:498:LEU:N	1:D:520:GLN:OE1	2.41	0.47
1:C:568:TYR:CD2	1:C:571:LYS:HD2	2.49	0.47
1:A:40:GLU:HG3	3:A:801:HOH:O	2.14	0.47
1:B:635:LEU:HD13	1:B:643:ILE:HG21	1.97	0.47
1:D:130:ARG:NH2	3:D:838:HOH:O	2.48	0.47
1:B:442:LEU:HD12	1:B:443:PRO:HD2	1.97	0.47
1:B:316:TYR:CE2	1:B:352:LEU:HG	2.50	0.46
1:B:302:GLY:HA2	1:B:333:TYR:O	2.15	0.46
1:B:573:VAL:O	1:B:587:ALA:HA	2.16	0.46
1:B:116:LEU:HD11	1:B:131:TYR:HB3	1.97	0.46
1:D:338:ASN:HD22	1:D:378:GLN:NE2	2.09	0.46
1:C:59:PRO:HD2	1:C:150:GLY:HA3	1.97	0.46
1:D:278:VAL:HG11	1:D:564:PHE:CD1	2.51	0.46
1:A:615:ASP:OD2	1:A:619:GLU:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:139:LYS:NZ	3:C:815:HOH:O	2.32	0.46
1:D:615:ASP:OD2	1:D:619:GLU:HB2	2.15	0.46
1:A:44:CYS:SG	1:A:57:LYS:HE3	2.56	0.46
1:B:485:ILE:O	1:B:488:THR:HG22	2.16	0.46
1:A:231:VAL:HG23	3:A:907:HOH:O	2.15	0.45
1:D:582:THR:HG22	3:D:802:HOH:O	2.16	0.45
1:D:519:PHE:O	1:D:523:THR:HG23	2.16	0.45
1:C:371:GLY:HA2	1:C:382:ASN:OD1	2.16	0.45
1:D:357:ARG:NH1	3:D:836:HOH:O	2.47	0.45
1:D:485:ILE:O	1:D:488:THR:HG22	2.16	0.45
1:C:573:VAL:O	1:C:587:ALA:HA	2.17	0.45
1:D:612:TYR:HB2	1:D:644:LEU:HB3	1.98	0.45
1:C:219:VAL:HB	1:C:444:ARG:HD3	1.98	0.44
1:C:612:TYR:HB2	1:C:644:LEU:HB3	1.98	0.44
1:C:86:VAL:HG12	1:C:88:GLN:HE22	1.82	0.44
1:D:602:GLU:HG2	1:D:603:MET:N	2.32	0.44
1:D:367:TYR:O	1:D:407:VAL:HA	2.17	0.44
1:D:216:GLU:OE1	1:D:266:ARG:NH1	2.51	0.44
1:D:425:LEU:HD13	1:D:450:TYR:CG	2.53	0.44
1:C:58:SER:HA	1:C:59:PRO:HD3	1.84	0.44
1:A:568:TYR:CD2	1:A:571:LYS:HD2	2.53	0.44
1:B:556:PRO:HG2	1:B:559:TRP:CH2	2.53	0.44
1:B:567:GLY:HA2	1:B:572:TYR:CE2	2.53	0.44
1:C:412:GLY:HA2	3:C:808:HOH:O	2.16	0.44
1:D:573:VAL:O	1:D:587:ALA:HA	2.17	0.44
1:A:66:ILE:O	1:A:114:GLN:NE2	2.51	0.44
1:B:406:LYS:HE3	1:B:438:HIS:CD2	2.53	0.44
1:B:612:TYR:HB2	1:B:644:LEU:HB3	1.99	0.43
1:D:42:LYS:HB2	1:D:42:LYS:HE3	1.74	0.43
1:B:628:LYS:HG3	1:B:629:GLU:N	2.33	0.43
1:D:396:TRP:HE3	1:D:397:LEU:HD12	1.81	0.43
1:B:603:MET:HG2	1:B:603:MET:H	1.59	0.43
1:D:592:LYS:HG2	1:D:592:LYS:O	2.19	0.43
1:A:342:THR:HG23	1:A:343:ARG:HG3	2.00	0.43
1:C:601:LEU:HD12	1:C:633:VAL:HG21	2.00	0.43
1:A:216:GLU:OE1	1:A:266:ARG:NH1	2.52	0.43
1:B:167:ALA:HA	1:B:176:PRO:HD3	2.01	0.43
1:A:442:LEU:HD12	1:A:443:PRO:HD2	2.01	0.43
1:B:367:TYR:O	1:B:407:VAL:HA	2.19	0.43
1:C:460:LEU:HD13	1:C:471:PHE:CE1	2.54	0.43
1:D:134:PRO:HA	1:D:260:PRO:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:82:LYS:HD2	1:D:100:ARG:NH2	2.33	0.43
1:D:559:TRP:CE2	1:D:577:ARG:HD3	2.54	0.43
1:D:369:SER:HB3	1:D:407:VAL:HG12	2.01	0.43
1:B:225:GLY:O	1:B:242:PRO:HD3	2.19	0.43
1:C:37:VAL:HG11	1:C:108:LEU:HD22	2.01	0.43
1:B:370:SER:HB3	1:B:374:ASN:HD21	1.83	0.42
1:C:337:ASP:HA	1:C:366:TRP:HB2	2.00	0.42
1:D:119:ILE:O	1:D:129:PHE:HA	2.19	0.42
1:C:139:LYS:HA	1:D:375:ASP:OD1	2.19	0.42
1:D:307:ILE:HG21	1:D:535:ALA:HA	2.01	0.42
1:D:205:ARG:NH1	1:D:207:GLY:O	2.52	0.42
1:D:302:GLY:HA2	1:D:333:TYR:O	2.19	0.42
1:D:635:LEU:HD13	1:D:643:ILE:HD13	2.00	0.42
1:A:305:SER:HA	1:A:322:TYR:CZ	2.54	0.42
1:A:191:ARG:CG	1:A:191:ARG:HH11	2.32	0.42
1:A:567:GLY:HA2	1:A:572:TYR:CE2	2.55	0.42
1:B:241:PHE:HB3	1:B:242:PRO:HD2	2.01	0.42
1:C:615:ASP:OD2	1:C:619:GLU:HB2	2.19	0.42
1:B:130:ARG:NH1	3:B:812:HOH:O	2.31	0.42
1:C:603:MET:HG2	1:C:603:MET:H	1.69	0.42
1:D:371:GLY:HA2	1:D:382:ASN:OD1	2.20	0.42
1:B:119:ILE:O	1:B:129:PHE:HA	2.20	0.42
1:C:213:LEU:HB3	1:C:269:THR:HB	2.02	0.42
1:C:375:ASP:HA	1:D:140:ALA:HB3	2.02	0.42
1:D:441:THR:OG1	1:D:444:ARG:NH1	2.46	0.42
1:C:567:GLY:HA2	1:C:572:TYR:CE2	2.55	0.42
1:D:403:LYS:HA	1:D:403:LYS:HD2	1.81	0.42
1:C:299:PHE:O	3:C:802:HOH:O	2.22	0.41
1:D:165:SER:HA	1:D:194:TYR:CD1	2.54	0.41
1:D:305:SER:HA	1:D:322:TYR:CZ	2.54	0.41
1:A:485:ILE:O	1:A:488:THR:HG22	2.20	0.41
1:A:464:ASN:HB3	1:A:471:PHE:CD2	2.55	0.41
1:B:554:ARG:NH2	3:B:838:HOH:O	2.49	0.41
1:C:269:THR:HG22	1:C:277:ILE:HD13	2.03	0.41
1:C:568:TYR:HA	1:C:569:PRO:HD3	1.88	0.41
1:C:608:THR:CG2	1:C:626:LYS:HG3	2.50	0.41
1:D:241:PHE:HB3	1:D:242:PRO:HD2	2.02	0.41
1:D:567:GLY:HA2	1:D:572:TYR:CE2	2.55	0.41
1:C:73:LEU:HD23	1:C:108:LEU:HD13	2.02	0.41
1:D:499:LEU:HD13	1:D:544:VAL:HG11	2.03	0.41
1:A:16:HIS:HB3	1:A:17:MET:H	1.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:367:TYR:OH	1:A:397:LEU:HD11	2.21	0.41
1:A:201:PRO:HB2	1:A:213:LEU:HD11	2.02	0.41
1:A:455:GLY:HA3	1:A:529:ASN:HD21	1.86	0.41
1:B:42:LYS:HE3	1:B:42:LYS:HB2	1.73	0.41
1:A:367:TYR:CD2	1:A:380:PRO:HG2	2.56	0.41
1:C:352:LEU:HA	1:C:352:LEU:HD23	1.86	0.41
1:D:583:TRP:CE3	1:D:603:MET:HE3	2.56	0.41
1:B:441:THR:OG1	1:B:444:ARG:NH1	2.43	0.41
1:C:291:TYR:CE2	1:C:451:PRO:HD2	2.56	0.41
1:A:119:ILE:O	1:A:129:PHE:HA	2.21	0.41
1:A:572:TYR:CD2	1:A:595:LEU:HD21	2.56	0.41
1:A:615:ASP:OD1	1:A:619:GLU:N	2.50	0.41
1:C:369:SER:HB3	1:C:407:VAL:HG12	2.03	0.41
1:D:378:GLN:HG3	1:D:378:GLN:H	1.69	0.41
1:A:291:TYR:CE2	1:A:451:PRO:HD2	2.56	0.40
1:D:367:TYR:CD2	1:D:380:PRO:HG2	2.56	0.40
1:C:332:GLU:HG2	3:C:802:HOH:O	2.20	0.40
1:A:248:ASN:HB2	1:A:373:TRP:O	2.21	0.40
1:C:574:VAL:HG13	1:C:587:ALA:HB2	2.03	0.40
1:D:264:PRO:HB2	1:D:266:ARG:HH12	1.86	0.40
1:D:66:ILE:O	1:D:114:GLN:NE2	2.55	0.40
1:A:302:GLY:HA2	1:A:333:TYR:O	2.21	0.40
1:A:241:PHE:HB3	1:A:242:PRO:HD2	2.04	0.40
1:A:147:GLU:CD	1:A:266:ARG:HH21	2.24	0.40
1:A:573:VAL:O	1:A:587:ALA:HA	2.21	0.40
1:B:351:SER:HA	1:B:354:GLU:OE1	2.22	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	630/642 (98%)	600 (95%)	29 (5%)	1 (0%)	47	58
1	B	626/642 (98%)	601 (96%)	23 (4%)	2 (0%)	41	50
1	C	630/642 (98%)	602 (96%)	27 (4%)	1 (0%)	47	58
1	D	627/642 (98%)	600 (96%)	26 (4%)	1 (0%)	47	58
All	All	2513/2568 (98%)	2403 (96%)	105 (4%)	5 (0%)	47	58

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	605	ALA
1	A	485	ILE
1	B	485	ILE
1	C	485	ILE
1	D	485	ILE

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	548/558 (98%)	543 (99%)	5 (1%)	78	89
1	B	544/558 (98%)	542 (100%)	2 (0%)	91	96
1	C	548/558 (98%)	545 (100%)	3 (0%)	88	95
1	D	545/558 (98%)	537 (98%)	8 (2%)	65	79
All	All	2185/2232 (98%)	2167 (99%)	18 (1%)	81	91

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	270	VAL
1	A	337	ASP
1	A	378	GLN
1	A	597	LEU
1	A	632	LYS
1	B	191	ARG

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Mol	Chain	Res	Type
1	B	597	LEU
1	C	88	GLN
1	C	597	LEU
1	C	608	THR
1	D	38	GLN
1	D	39	ASN
1	D	82	LYS
1	D	597	LEU
1	D	600	ASP
1	D	603	MET
1	D	616	LYS
1	D	626	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	529	ASN
1	B	338	ASN
1	B	378	GLN
1	C	338	ASN
1	D	338	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	632/642 (98%)	0.26	15 (2%) 59 66	14, 24, 46, 68	0
1	B	628/642 (97%)	0.33	31 (4%) 29 36	17, 29, 50, 68	0
1	C	632/642 (98%)	0.36	29 (4%) 32 39	14, 25, 46, 67	0
1	D	629/642 (97%)	0.51	49 (7%) 13 17	18, 31, 56, 77	0
All	All	2521/2568 (98%)	0.36	124 (4%) 29 36	14, 27, 50, 77	0

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	605	ALA	6.2
1	D	600	ASP	5.7
1	A	605	ALA	5.5
1	B	605	ALA	5.3
1	D	618	GLY	5.0
1	C	605	ALA	4.8
1	D	191	ARG	4.6
1	D	626	LYS	4.5
1	D	635	LEU	4.4
1	B	594	PRO	4.1
1	A	627	VAL	4.0
1	B	616	LYS	4.0
1	B	633	VAL	4.0
1	D	634	GLN	3.9
1	A	16	HIS	3.9
1	A	603	MET	3.9
1	B	610	ALA	3.9
1	A	628	LYS	3.9
1	B	628	LYS	3.8
1	B	632	LYS	3.7
1	D	617	LYS	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	629	GLU	3.6
1	C	626	LYS	3.6
1	A	604	PHE	3.6
1	B	617	LYS	3.5
1	D	627	VAL	3.5
1	A	600	ASP	3.5
1	C	592	LYS	3.4
1	C	527	LEU	3.4
1	D	39	ASN	3.4
1	C	608	THR	3.4
1	C	635	LEU	3.4
1	D	19	LEU	3.4
1	B	646	ILE	3.3
1	B	631	GLY	3.3
1	D	637	ILE	3.3
1	C	628	LYS	3.3
1	D	603	MET	3.2
1	D	630	ASN	3.2
1	D	595	LEU	3.2
1	D	628	LYS	3.2
1	C	625	LEU	3.2
1	C	347	GLN	3.1
1	C	593	GLU	3.1
1	A	85	THR	3.1
1	D	347	GLN	3.0
1	D	606	GLY	3.0
1	D	616	LYS	3.0
1	D	604	PHE	3.0
1	D	421	TYR	3.0
1	D	597	LEU	3.0
1	D	155	GLU	3.0
1	D	598	LYS	2.9
1	C	351	SER	2.9
1	B	581	ASP	2.9
1	D	608	THR	2.9
1	D	611	LEU	2.9
1	C	358	ASP	2.9
1	D	625	LEU	2.9
1	D	68	ASP	2.8
1	B	599	LEU	2.8
1	D	607	LYS	2.8
1	D	624	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	614	ASP	2.8
1	B	598	LYS	2.8
1	C	16	HIS	2.8
1	C	624	SER	2.8
1	C	354	GLU	2.8
1	A	420	LEU	2.7
1	C	602	GLU	2.7
1	B	39	ASN	2.7
1	B	600	ASP	2.7
1	B	604	PHE	2.7
1	A	630	ASN	2.7
1	C	631	GLY	2.7
1	C	632	LYS	2.7
1	B	595	LEU	2.7
1	D	633	VAL	2.7
1	D	407	VAL	2.6
1	D	620	PRO	2.6
1	D	636	GLU	2.6
1	D	417	THR	2.5
1	B	629	GLU	2.5
1	D	583	TRP	2.5
1	B	634	GLN	2.5
1	A	633	VAL	2.5
1	C	599	LEU	2.4
1	A	626	LYS	2.4
1	B	647	LYS	2.4
1	B	626	LYS	2.4
1	C	598	LYS	2.4
1	C	594	PRO	2.3
1	D	599	LEU	2.3
1	C	609	VAL	2.3
1	D	638	ARG	2.3
1	D	446	TRP	2.3
1	B	565	VAL	2.3
1	C	606	GLY	2.3
1	B	321	ARG	2.3
1	B	592	LYS	2.3
1	D	348	ARG	2.3
1	C	636	GLU	2.2
1	B	645	CYS	2.2
1	A	424	ILE	2.2
1	A	629	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	350	LYS	2.2
1	B	518	VAL	2.2
1	D	187	PRO	2.2
1	D	592	LYS	2.2
1	B	345	GLY	2.2
1	C	630	ASN	2.2
1	B	596	LYS	2.1
1	A	616	LYS	2.1
1	D	455	GLY	2.1
1	B	348	ARG	2.1
1	D	632	LYS	2.1
1	D	615	ASP	2.1
1	C	554	ARG	2.1
1	D	170	GLY	2.0
1	D	456	SER	2.0
1	D	629	GLU	2.0
1	D	157	THR	2.0
1	C	627	VAL	2.0
1	B	630	ASN	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 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(\AA^2)	Q<0.9
2	CA	B	701	1/1	0.97	0.05	27,27,27,27	0
2	CA	C	701	1/1	0.97	0.10	21,21,21,21	0
2	CA	D	701	1/1	0.98	0.08	25,25,25,25	0
2	CA	A	701	1/1	0.99	0.07	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.