



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:21 PM UTC

PDB ID : 3CZL / pdb_00003czl
Title : Crystal Structure Analysis of Sucrose hydrolase(SUH) E322Q-glucose complex
Authors : Kim, M.I.; Rhee, S.
Deposited on : 2008-04-29
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 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

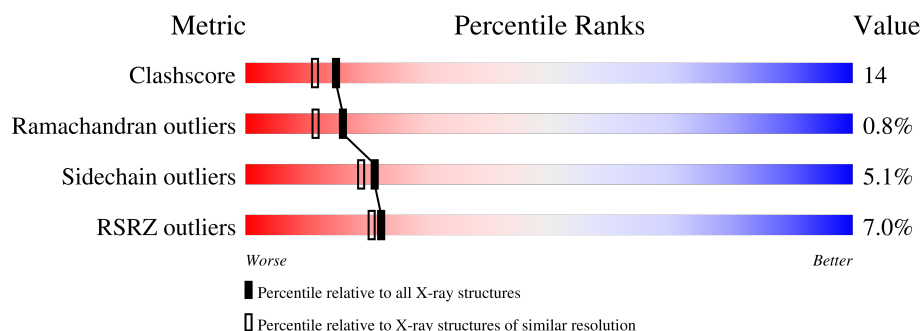
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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	644	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5162 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.

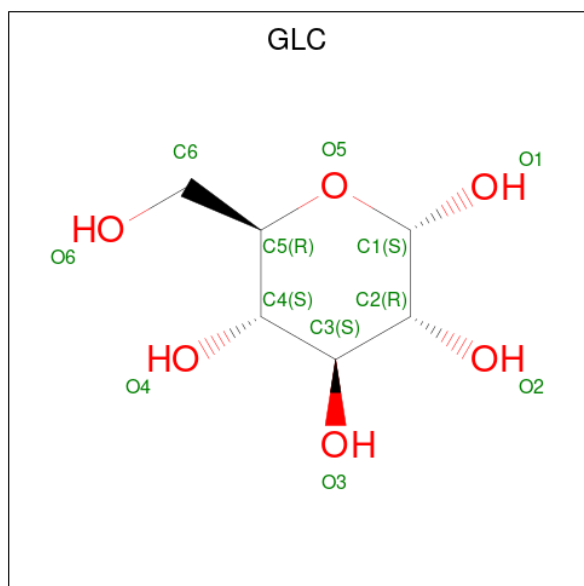
- Molecule 1 is a protein called sucrose hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	612	4752	3015	858	860	19	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	322	GLN	GLU	engineered mutation	UNP Q6UVM5

- Molecule 2 is alpha-D-glucopyranose (CCD ID: GLC) (formula: C₆H₁₂O₆).



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

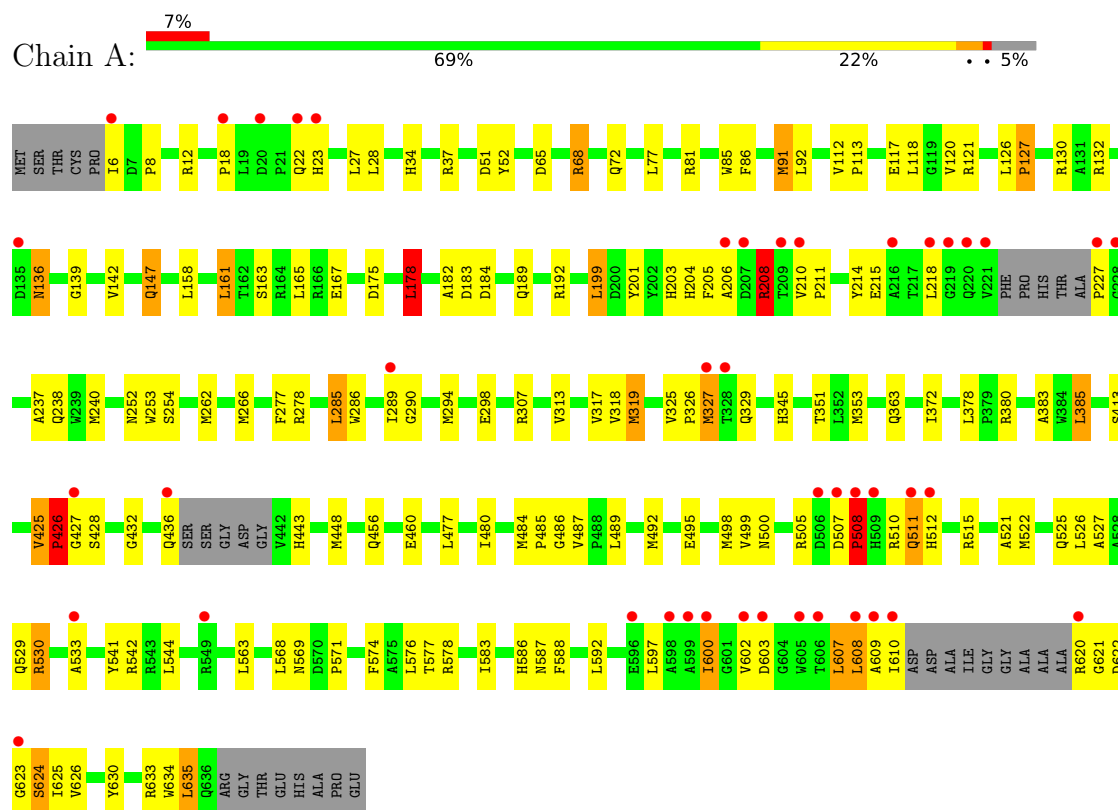
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	398	Total 398	O 398	0	0

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: sucrose hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	96.80Å 119.10Å 55.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 50.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.0 (50.00-2.00) 97.8 (50.00-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.42 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.203 , 0.248 0.210 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.7	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	0.94	EDS
Total number of atoms	5162	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.18% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.73	21/4881 (0.4%)	1.02	25/6661 (0.4%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	266	MET	SD-CE	7.19	1.97	1.79
1	A	353	MET	SD-CE	6.79	1.96	1.79
1	A	327	MET	CG-SD	6.77	1.97	1.80
1	A	484	MET	SD-CE	6.75	1.96	1.79
1	A	327	MET	SD-CE	6.63	1.96	1.79
1	A	492	MET	SD-CE	6.50	1.95	1.79
1	A	240	MET	SD-CE	6.39	1.95	1.79
1	A	522	MET	CG-SD	6.30	1.96	1.80
1	A	498	MET	SD-CE	6.29	1.95	1.79
1	A	262	MET	SD-CE	6.28	1.95	1.79
1	A	91	MET	SD-CE	6.09	1.94	1.79
1	A	319	MET	SD-CE	6.00	1.94	1.79
1	A	522	MET	SD-CE	5.98	1.94	1.79
1	A	353	MET	CG-SD	5.95	1.95	1.80
1	A	294	MET	SD-CE	5.89	1.94	1.79
1	A	262	MET	CG-SD	5.83	1.95	1.80
1	A	448	MET	SD-CE	5.82	1.94	1.79
1	A	448	MET	CG-SD	5.53	1.94	1.80
1	A	484	MET	CG-SD	5.20	1.93	1.80
1	A	294	MET	CG-SD	5.16	1.93	1.80
1	A	492	MET	CG-SD	5.04	1.93	1.80

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	VAL	N-CA-C	8.47	120.56	111.58
1	A	428	SER	N-CA-C	8.00	122.02	110.10
1	A	285	LEU	N-CA-C	7.79	120.75	111.33
1	A	609	ALA	N-CA-C	-7.67	103.38	112.89
1	A	511	GLN	N-CA-C	-7.60	100.53	110.33
1	A	199	LEU	N-CA-C	-7.42	103.27	111.36
1	A	385	LEU	N-CA-C	-7.01	96.21	108.20
1	A	52	TYR	N-CA-C	6.88	118.44	111.07
1	A	525	GLN	N-CA-C	-6.45	103.54	111.40
1	A	8	PRO	N-CA-C	6.16	118.22	110.70
1	A	622	ASP	N-CA-C	-6.00	104.41	113.89
1	A	201	TYR	N-CA-C	-5.96	105.78	113.16
1	A	487	VAL	N-CA-C	5.62	113.62	107.60
1	A	427	GLY	N-CA-C	5.57	124.74	115.66
1	A	425	VAL	CA-C-N	5.56	126.79	119.84
1	A	425	VAL	C-N-CA	5.56	126.79	119.84
1	A	254	SER	N-CA-C	-5.49	106.53	113.18
1	A	432	GLY	N-CA-C	-5.49	105.05	112.57
1	A	178	LEU	N-CA-C	5.40	119.71	113.18
1	A	484	MET	N-CA-C	5.36	117.92	110.20
1	A	521	ALA	N-CA-C	-5.29	101.84	110.20
1	A	492	MET	N-CA-C	5.28	118.45	110.64
1	A	142	VAL	N-CA-C	5.20	115.75	108.89
1	A	208	ARG	N-CA-C	-5.10	106.90	113.02
1	A	413	SER	N-CA-C	5.09	116.90	108.20

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	4752	0	4575	129	0
2	A	12	0	12	0	0
3	A	398	0	0	12	0
All	All	5162	0	4587	129	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 14.

All (129) 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:208:ARG:HH11	1:A:208:ARG:HB3	1.10	1.09
1:A:563:LEU:HB2	1:A:578:ARG:HH11	1.34	0.90
1:A:507:ASP:HB2	1:A:510:ARG:HH21	1.39	0.86
1:A:533:ALA:HB1	1:A:542:ARG:HH21	1.44	0.82
1:A:319:MET:H	1:A:345:HIS:HD2	1.24	0.82
1:A:208:ARG:HH11	1:A:208:ARG:CB	1.91	0.80
1:A:158:LEU:HD23	1:A:158:LEU:O	1.82	0.80
1:A:68:ARG:HG3	1:A:68:ARG:HH11	1.46	0.79
1:A:597:LEU:HD22	1:A:602:VAL:HG23	1.63	0.79
1:A:456:GLN:HE21	1:A:499:VAL:HB	1.48	0.78
1:A:597:LEU:HA	1:A:600:ILE:HD11	1.64	0.78
1:A:563:LEU:HB2	1:A:578:ARG:NH1	1.99	0.76
1:A:130:ARG:HE	1:A:147:GLN:HG2	1.49	0.76
1:A:508:PRO:HA	3:A:1030:HOH:O	1.88	0.74
1:A:158:LEU:HD23	1:A:158:LEU:C	2.12	0.73
1:A:130:ARG:HH21	1:A:147:GLN:HE21	1.37	0.73
1:A:505:ARG:HA	1:A:511:GLN:HG3	1.69	0.72
1:A:208:ARG:HB3	1:A:208:ARG:NH1	1.96	0.71
1:A:515:ARG:HD3	3:A:1019:HOH:O	1.91	0.70
1:A:600:ILE:H	1:A:600:ILE:HD13	1.56	0.69
1:A:163:SER:O	1:A:167:GLU:HG3	1.96	0.65
1:A:65:ASP:HB2	3:A:965:HOH:O	1.96	0.65
1:A:178:LEU:HD13	1:A:285:LEU:HD21	1.79	0.65
1:A:204:HIS:HD2	1:A:238:GLN:OE1	1.80	0.64
1:A:511:GLN:O	1:A:512:HIS:HB2	1.96	0.64
1:A:507:ASP:CG	1:A:510:ARG:HE	2.07	0.62
1:A:130:ARG:HH21	1:A:147:GLN:NE2	1.97	0.61
1:A:326:PRO:CG	1:A:329:GLN:HE21	2.12	0.61
1:A:610:ILE:HG12	1:A:610:ILE:O	2.01	0.60
1:A:237:ALA:O	1:A:238:GLN:HG3	2.02	0.60
1:A:456:GLN:O	1:A:460:GLU:HG3	2.03	0.59
1:A:158:LEU:C	1:A:158:LEU:CD2	2.76	0.58
1:A:607:LEU:HD23	1:A:634:TRP:HB3	1.86	0.58
1:A:132:ARG:HD2	1:A:183:ASP:OD1	2.04	0.57
1:A:577:THR:HG22	1:A:583:ILE:HD13	1.86	0.57
1:A:507:ASP:CB	1:A:510:ARG:HE	2.17	0.57
1:A:597:LEU:HA	1:A:600:ILE:CD1	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:ARG:NH2	1:A:28:LEU:HD13	2.20	0.57
1:A:325:VAL:HB	1:A:326:PRO:HD2	1.87	0.57
1:A:6:ILE:HG13	3:A:983:HOH:O	2.05	0.56
1:A:425:VAL:HG23	1:A:426:PRO:HD2	1.88	0.56
1:A:289:ILE:HG12	1:A:290:GLY:N	2.22	0.55
1:A:68:ARG:HH11	1:A:68:ARG:CG	2.17	0.55
1:A:277:PHE:CE2	1:A:317:VAL:HG11	2.42	0.55
1:A:326:PRO:HG2	1:A:329:GLN:HE21	1.72	0.54
1:A:577:THR:HG22	1:A:583:ILE:CD1	2.38	0.54
1:A:363:GLN:HA	1:A:588:PHE:O	2.06	0.54
1:A:380:ARG:HD3	3:A:928:HOH:O	2.06	0.54
1:A:210:VAL:HB	1:A:211:PRO:HD3	1.90	0.54
1:A:319:MET:H	1:A:345:HIS:CD2	2.15	0.54
1:A:587:ASN:O	1:A:630:TYR:HA	2.08	0.54
1:A:6:ILE:HG22	1:A:6:ILE:O	2.09	0.53
1:A:533:ALA:CB	1:A:542:ARG:HH21	2.20	0.53
1:A:117:GLU:OE1	1:A:530:ARG:NH1	2.41	0.53
1:A:277:PHE:HE2	1:A:317:VAL:HG11	1.73	0.52
1:A:436:GLN:HB2	1:A:443:HIS:CG	2.43	0.52
1:A:205:PHE:CE1	1:A:210:VAL:HG11	2.44	0.52
1:A:23:HIS:NE2	1:A:72:GLN:OE1	2.40	0.52
1:A:189:GLN:OE1	1:A:192:ARG:HD2	2.09	0.51
1:A:569:ASN:HB2	3:A:1023:HOH:O	2.11	0.51
1:A:81:ARG:HD3	1:A:85:TRP:CE2	2.46	0.51
1:A:436:GLN:HB2	1:A:443:HIS:CD2	2.45	0.51
1:A:204:HIS:CD2	1:A:238:GLN:OE1	2.63	0.51
1:A:34:HIS:ND1	1:A:37:ARG:NH2	2.50	0.51
1:A:507:ASP:CB	1:A:510:ARG:HH21	2.17	0.51
1:A:206:ALA:HA	1:A:238:GLN:HG2	1.93	0.51
1:A:608:LEU:O	1:A:608:LEU:HD22	2.10	0.50
1:A:65:ASP:OD1	1:A:68:ARG:NH2	2.45	0.50
1:A:18:PRO:HB2	1:A:65:ASP:HA	1.94	0.50
1:A:161:LEU:HD22	1:A:165:LEU:HG	1.93	0.50
1:A:625:ILE:HD13	1:A:635:LEU:HD11	1.94	0.50
1:A:307:ARG:CZ	1:A:345:HIS:CE1	2.95	0.49
1:A:326:PRO:HD2	1:A:329:GLN:NE2	2.27	0.49
1:A:526:LEU:HD12	1:A:529:GLN:NE2	2.27	0.49
1:A:600:ILE:HD13	1:A:600:ILE:N	2.26	0.49
1:A:132:ARG:HG3	1:A:184:ASP:OD1	2.12	0.49
1:A:620:ARG:HG3	1:A:621:GLY:N	2.27	0.49
1:A:600:ILE:HG12	1:A:602:VAL:HG22	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:PRO:HB3	1:A:578:ARG:NH2	2.28	0.48
1:A:485:PRO:HD3	1:A:578:ARG:NH1	2.28	0.48
1:A:530:ARG:HD3	1:A:541:TYR:CE2	2.49	0.47
1:A:130:ARG:NH2	1:A:147:GLN:HE21	2.08	0.47
1:A:318:VAL:HA	1:A:345:HIS:CD2	2.50	0.47
1:A:327:MET:HG3	1:A:351:THR:HG21	1.96	0.47
1:A:178:LEU:HB2	1:A:253:TRP:HZ2	1.79	0.47
1:A:592:LEU:HD11	1:A:626:VAL:CG1	2.45	0.47
1:A:608:LEU:O	1:A:608:LEU:HD13	2.16	0.46
1:A:326:PRO:CG	1:A:329:GLN:NE2	2.79	0.46
1:A:385:LEU:HD11	1:A:489:LEU:CD2	2.46	0.46
1:A:211:PRO:O	1:A:215:GLU:HG2	2.16	0.45
1:A:132:ARG:HG3	1:A:182:ALA:HB1	1.98	0.45
1:A:286:TRP:CG	1:A:298:GLU:HB2	2.51	0.45
1:A:495:GLU:HA	1:A:544:LEU:HD12	1.98	0.45
1:A:86:PHE:HA	1:A:91:MET:HE2	1.99	0.45
1:A:136:ASN:HD22	1:A:139:GLY:H	1.63	0.45
1:A:529:GLN:HG2	3:A:1196:HOH:O	2.17	0.45
1:A:120:VAL:O	1:A:121:ARG:HD2	2.17	0.45
1:A:203:HIS:CE1	1:A:252:ASN:HA	2.52	0.44
1:A:489:LEU:HD21	3:A:982:HOH:O	2.18	0.44
1:A:345:HIS:HE1	3:A:823:HOH:O	2.00	0.44
1:A:456:GLN:HE22	1:A:500:ASN:HB2	1.82	0.44
1:A:485:PRO:HB3	1:A:578:ARG:HH22	1.83	0.44
1:A:378:LEU:HD13	1:A:383:ALA:HA	2.00	0.44
1:A:86:PHE:CZ	1:A:345:HIS:HB3	2.53	0.43
1:A:425:VAL:CG2	1:A:426:PRO:HD2	2.48	0.43
1:A:175:ASP:OD1	1:A:278:ARG:NH1	2.43	0.43
1:A:526:LEU:HD12	1:A:529:GLN:HE22	1.83	0.43
1:A:385:LEU:HD11	1:A:489:LEU:HD22	2.00	0.43
1:A:608:LEU:H	1:A:608:LEU:HD12	1.84	0.43
1:A:18:PRO:O	1:A:68:ARG:HG3	2.19	0.42
1:A:214:TYR:O	1:A:218:LEU:HG	2.19	0.42
1:A:586:HIS:HE1	3:A:848:HOH:O	2.03	0.42
1:A:383:ALA:HB1	1:A:486:GLY:HA2	2.01	0.42
1:A:383:ALA:HB1	1:A:486:GLY:CA	2.50	0.42
1:A:620:ARG:CG	1:A:621:GLY:N	2.83	0.41
1:A:51:ASP:C	1:A:51:ASP:OD1	2.63	0.41
1:A:68:ARG:CG	1:A:68:ARG:NH1	2.81	0.41
1:A:623:GLY:O	1:A:624:SER:C	2.63	0.41
1:A:126:LEU:HB3	1:A:127:PRO:HD2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:VAL:N	1:A:113:PRO:CD	2.83	0.41
1:A:485:PRO:HB3	1:A:578:ARG:NH1	2.35	0.41
1:A:27:LEU:HD23	3:A:1010:HOH:O	2.20	0.40
1:A:132:ARG:NH2	3:A:934:HOH:O	2.52	0.40
1:A:136:ASN:HD22	1:A:136:ASN:C	2.28	0.40
1:A:372:ILE:HG13	1:A:574:PHE:HZ	1.85	0.40
1:A:507:ASP:HB3	1:A:510:ARG:HE	1.85	0.40
1:A:511:GLN:O	1:A:512:HIS:CB	2.61	0.40
1:A:527:ALA:O	1:A:530:ARG:HB3	2.21	0.40
1:A:569:ASN:O	1:A:571:PRO:HD3	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	604/644 (94%)	571 (94%)	28 (5%)	5 (1%)	16 11

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	508	PRO
1	A	603	ASP
1	A	624	SER
1	A	127	PRO
1	A	426	PRO

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	470/491 (96%)	446 (95%)	24 (5%)	21	19

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	68	ARG
1	A	77	LEU
1	A	92	LEU
1	A	118	LEU
1	A	136	ASN
1	A	147	GLN
1	A	161	LEU
1	A	178	LEU
1	A	199	LEU
1	A	208	ARG
1	A	227	PRO
1	A	426	PRO
1	A	477	LEU
1	A	480	ILE
1	A	508	PRO
1	A	530	ARG
1	A	568	LEU
1	A	576	LEU
1	A	600	ILE
1	A	607	LEU
1	A	608	LEU
1	A	633	ARG
1	A	635	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	69	GLN
1	A	88	GLN
1	A	106	GLN
1	A	136	ASN

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Mol	Chain	Res	Type
1	A	147	GLN
1	A	203	HIS
1	A	204	HIS
1	A	295	ASN
1	A	329	GLN
1	A	345	HIS
1	A	401	HIS
1	A	456	GLN
1	A	459	GLN
1	A	511	GLN
1	A	529	GLN
1	A	531	HIS
1	A	550	GLN
1	A	586	HIS
1	A	587	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	GLC	A	3497	-	12,12,12	1.05	0	17,17,17	0.82	1 (5%)

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	GLC	A	3497	-	-	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	3497	GLC	C6-C5-C4	-2.02	108.06	113.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	612/644 (95%)	0.40	43 (7%) 22 21	15, 25, 46, 56	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	605	TRP	5.6
1	A	219	GLY	4.9
1	A	506	ASP	4.2
1	A	599	ALA	4.1
1	A	609	ALA	4.1
1	A	289	ILE	3.9
1	A	610	ILE	3.7
1	A	227	PRO	3.5
1	A	207	ASP	3.4
1	A	603	ASP	3.4
1	A	623	GLY	3.4
1	A	436	GLN	3.2
1	A	511	GLN	3.1
1	A	20	ASP	2.9
1	A	600	ILE	2.9
1	A	220	GLN	2.9
1	A	228	GLY	2.9
1	A	6	ILE	2.7
1	A	512	HIS	2.7
1	A	221	VAL	2.7
1	A	206	ALA	2.5
1	A	23	HIS	2.5
1	A	598	ALA	2.5
1	A	508	PRO	2.5
1	A	509	HIS	2.5
1	A	608	LEU	2.4
1	A	22	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	209	THR	2.4
1	A	327	MET	2.4
1	A	18	PRO	2.3
1	A	596	GLU	2.3
1	A	135	ASP	2.3
1	A	620	ARG	2.3
1	A	602	VAL	2.3
1	A	328	THR	2.3
1	A	507	ASP	2.3
1	A	427	GLY	2.2
1	A	210	VAL	2.1
1	A	216	ALA	2.1
1	A	533	ALA	2.1
1	A	218	LEU	2.0
1	A	549	ARG	2.0
1	A	606	THR	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
2	GLC	A	3497	12/12	0.80	0.14	23,28,32,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.