



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:33 PM UTC

PDB ID : 3CZK / pdb_00003czk
Title : Crystal Structure Analysis of Sucrose hydrolase(SUH) E322Q-sucrose complex
Authors : Kim, M.I.; Rhee, S.
Deposited on : 2008-04-29
Resolution : 2.20 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

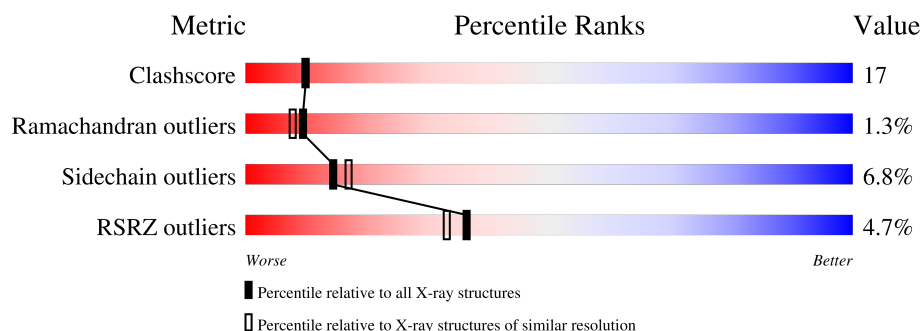
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	5% 62% 29% 5% .
2	B	2	100%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5085 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.

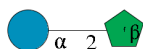
- Molecule 1 is a protein called Sucrose hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	618	Total	C	N	O	S	0	0	0
			4798	3043	869	867	19			

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 an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	2	Total	C	O	0	0	0
			23	12	11			

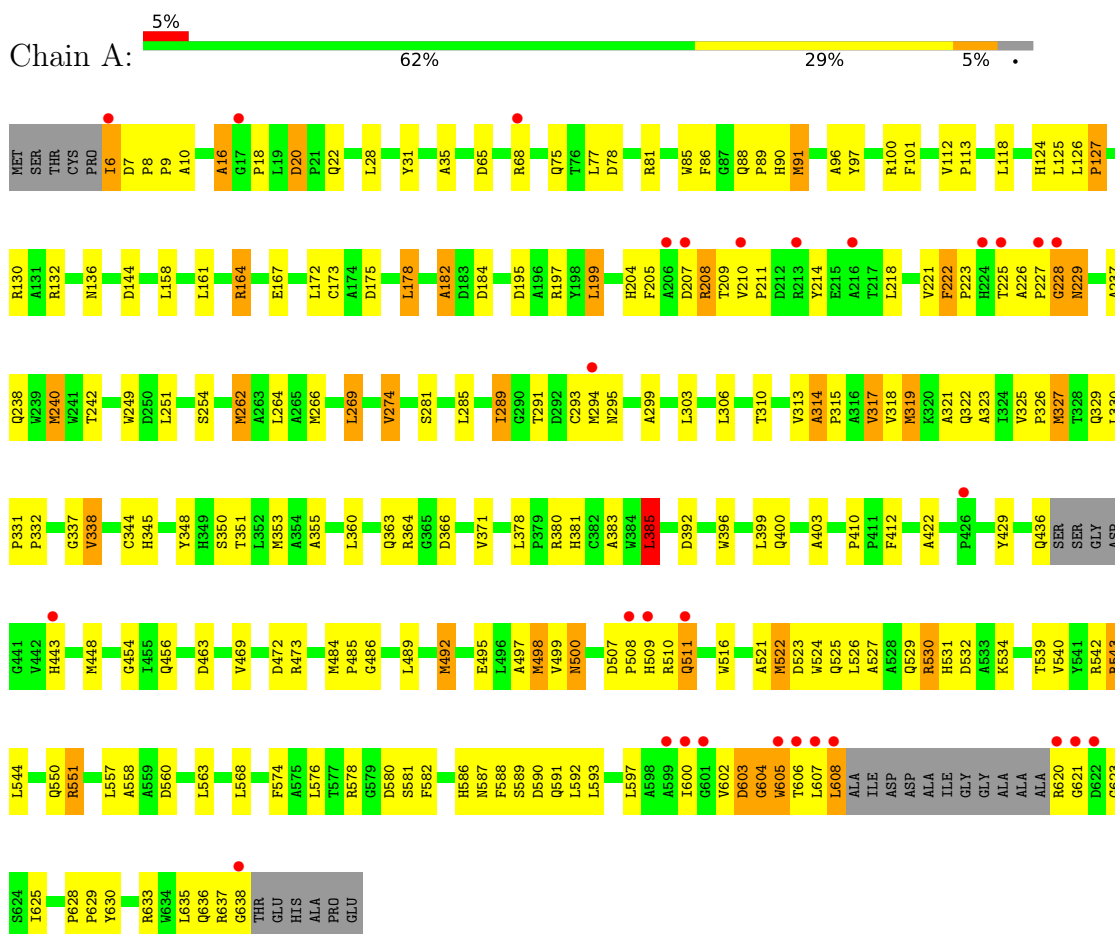
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	264	Total	O	0	0
			264	264		

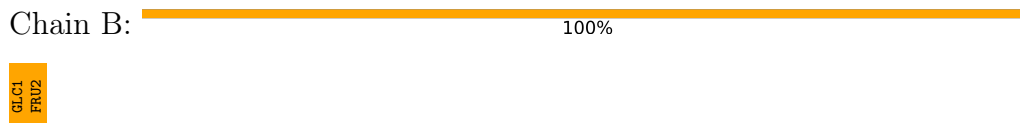
3 Residue-property plots [i](#)

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



• Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	96.50Å 118.60Å 55.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	83.5 (50.00-2.20) 94.3 (50.00-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.200 , 0.262 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.635	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5085	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% 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: FRU, 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	0.62	22/4931 (0.4%)	0.96	22/6731 (0.3%)

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	91	MET	SD-CE	6.45	1.95	1.79
1	A	319	MET	SD-CE	6.35	1.95	1.79
1	A	353	MET	SD-CE	6.19	1.95	1.79
1	A	492	MET	SD-CE	6.19	1.95	1.79
1	A	240	MET	SD-CE	6.18	1.95	1.79
1	A	448	MET	SD-CE	6.09	1.94	1.79
1	A	262	MET	SD-CE	6.06	1.94	1.79
1	A	353	MET	CG-SD	6.03	1.95	1.80
1	A	294	MET	SD-CE	5.97	1.94	1.79
1	A	266	MET	SD-CE	5.93	1.94	1.79
1	A	484	MET	SD-CE	5.91	1.94	1.79
1	A	498	MET	SD-CE	5.84	1.94	1.79
1	A	327	MET	SD-CE	5.81	1.94	1.79
1	A	522	MET	CG-SD	5.75	1.95	1.80
1	A	262	MET	CG-SD	5.75	1.95	1.80
1	A	448	MET	CG-SD	5.61	1.94	1.80
1	A	522	MET	SD-CE	5.41	1.93	1.79
1	A	492	MET	CG-SD	5.32	1.94	1.80
1	A	91	MET	CG-SD	5.31	1.94	1.80
1	A	240	MET	CG-SD	5.23	1.93	1.80
1	A	498	MET	CG-SD	5.14	1.93	1.80
1	A	484	MET	CG-SD	5.08	1.93	1.80

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	VAL	N-CA-C	7.56	119.59	111.58
1	A	16	ALA	N-CA-C	-7.33	101.10	110.19
1	A	199	LEU	N-CA-C	-7.18	103.49	111.82
1	A	385	LEU	N-CA-C	-6.97	96.27	108.20
1	A	500	ASN	N-CA-C	6.68	118.95	110.33
1	A	314	ALA	CA-C-N	6.66	126.35	119.56
1	A	314	ALA	C-N-CA	6.66	126.35	119.56
1	A	229	ASN	N-CA-C	-6.13	105.32	114.64
1	A	208	ARG	N-CA-C	-6.06	105.53	113.17
1	A	525	GLN	N-CA-C	-5.91	104.79	112.23
1	A	605	TRP	N-CA-C	5.70	122.94	110.80
1	A	396	TRP	N-CA-C	5.70	117.49	111.28
1	A	422	ALA	N-CA-C	-5.56	106.50	113.28
1	A	604	GLY	N-CA-C	5.55	124.19	111.04
1	A	8	PRO	N-CA-C	5.47	117.38	110.70
1	A	182	ALA	N-CA-C	-5.22	103.13	110.50
1	A	228	GLY	CA-C-N	5.17	130.88	121.41
1	A	228	GLY	C-N-CA	5.17	130.88	121.41
1	A	254	SER	N-CA-C	-5.09	107.07	113.28
1	A	222	PHE	CA-C-N	5.03	125.30	119.47
1	A	222	PHE	C-N-CA	5.03	125.30	119.47
1	A	492	MET	N-CA-C	5.01	118.06	110.64

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	4798	0	4613	164	0
2	B	23	0	21	2	0
3	A	264	0	0	9	1
All	All	5085	0	4634	165	1

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 17.

All (165) 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:550:GLN:HE21	1:A:608:LEU:HB2	1.17	1.08
1:A:18:PRO:O	1:A:68:ARG:HD2	1.56	1.05
1:A:226:ALA:C	1:A:228:GLY:H	1.66	0.97
1:A:456:GLN:HE21	1:A:499:VAL:HB	1.30	0.96
1:A:281:SER:H	1:A:322:GLN:NE2	1.73	0.87
1:A:96:ALA:HB2	1:A:492:MET:HB2	1.63	0.81
1:A:498:MET:HE3	1:A:523:ASP:HB2	1.66	0.78
1:A:605:TRP:HE3	1:A:606:THR:H	1.31	0.78
1:A:319:MET:H	1:A:345:HIS:HD2	1.32	0.77
1:A:550:GLN:NE2	1:A:608:LEU:HB2	1.97	0.77
1:A:178:LEU:HD13	1:A:285:LEU:HD21	1.66	0.76
1:A:299:ALA:O	1:A:303:LEU:HD13	1.86	0.75
1:A:226:ALA:C	1:A:228:GLY:N	2.37	0.75
1:A:226:ALA:O	1:A:228:GLY:N	2.20	0.74
1:A:602:VAL:O	1:A:603:ASP:HB2	1.90	0.72
1:A:563:LEU:HD11	1:A:576:LEU:HD21	1.71	0.70
1:A:16:ALA:HB2	1:A:28:LEU:HD21	1.73	0.70
1:A:214:TYR:O	1:A:218:LEU:HG	1.91	0.69
1:A:403:ALA:HA	1:A:412:PHE:CE1	2.29	0.68
1:A:132:ARG:HG3	1:A:184:ASP:CG	2.20	0.66
1:A:497:ALA:HB2	1:A:540:VAL:HG21	1.78	0.65
1:A:392:ASP:OD2	2:B:1:GLC:O2	2.10	0.65
1:A:225:THR:HB	1:A:436:GLN:HE22	1.61	0.64
1:A:89:PRO:O	1:A:551:ARG:HD3	1.99	0.62
1:A:542:ARG:HG2	1:A:542:ARG:HH11	1.63	0.62
1:A:318:VAL:HG13	1:A:345:HIS:HB2	1.81	0.62
1:A:485:PRO:HB3	1:A:578:ARG:NH1	2.14	0.61
1:A:485:PRO:HB3	1:A:578:ARG:HH12	1.65	0.61
1:A:97:TYR:HB2	1:A:100:ARG:HG2	1.82	0.61
1:A:330:LEU:HB2	1:A:331:PRO:HD3	1.82	0.60
1:A:563:LEU:CD1	1:A:576:LEU:HD21	2.32	0.60
1:A:597:LEU:HD11	1:A:623:GLY:H	1.67	0.59
1:A:68:ARG:HD3	3:A:970:HOH:O	2.03	0.59
1:A:325:VAL:HB	1:A:326:PRO:HD2	1.84	0.59
1:A:319:MET:H	1:A:345:HIS:CD2	2.19	0.58
1:A:269:LEU:O	1:A:274:VAL:HG13	2.03	0.58
1:A:550:GLN:NE2	1:A:608:LEU:H	2.00	0.58
1:A:16:ALA:O	1:A:18:PRO:HD3	2.04	0.57
1:A:592:LEU:HD12	1:A:628:PRO:HA	1.86	0.57
1:A:600:ILE:HG13	1:A:602:VAL:HG22	1.87	0.57
1:A:65:ASP:HB2	3:A:853:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:625:ILE:HD13	1:A:635:LEU:HD11	1.88	0.56
1:A:281:SER:H	1:A:322:GLN:HE21	1.53	0.56
1:A:606:THR:HG22	1:A:620:ARG:HB2	1.87	0.56
1:A:210:VAL:HB	1:A:211:PRO:HD3	1.88	0.56
1:A:158:LEU:HD23	1:A:158:LEU:C	2.31	0.56
1:A:295:ASN:HD21	1:A:325:VAL:HG11	1.70	0.56
1:A:360:LEU:HD11	1:A:473:ARG:HD3	1.88	0.56
1:A:456:GLN:NE2	1:A:500:ASN:H	2.03	0.56
1:A:348:TYR:HA	1:A:385:LEU:HB3	1.88	0.55
1:A:508:PRO:HA	1:A:511:GLN:CD	2.32	0.54
1:A:337:GLY:HA3	3:A:859:HOH:O	2.06	0.54
1:A:550:GLN:HE21	1:A:608:LEU:H	1.55	0.54
1:A:225:THR:HG21	3:A:1025:HOH:O	2.07	0.54
1:A:269:LEU:HD22	3:A:819:HOH:O	2.07	0.54
1:A:380:ARG:HG3	1:A:381:HIS:CD2	2.42	0.54
1:A:558:ALA:O	1:A:578:ARG:NH2	2.41	0.54
1:A:281:SER:H	1:A:322:GLN:HE22	1.54	0.53
1:A:175:ASP:HB2	3:A:806:HOH:O	2.08	0.52
1:A:325:VAL:HG21	1:A:329:GLN:HB3	1.90	0.52
1:A:587:ASN:O	1:A:630:TYR:HA	2.10	0.52
1:A:378:LEU:HD13	1:A:383:ALA:HA	1.93	0.51
1:A:587:ASN:ND2	1:A:589:SER:H	2.09	0.51
1:A:606:THR:HA	1:A:635:LEU:HD23	1.91	0.51
1:A:158:LEU:HD23	1:A:158:LEU:O	2.09	0.51
1:A:604:GLY:HA3	1:A:638:GLY:C	2.35	0.51
1:A:221:VAL:C	1:A:223:PRO:HD3	2.36	0.50
1:A:542:ARG:HG2	1:A:542:ARG:NH1	2.26	0.50
1:A:130:ARG:HB2	1:A:144:ASP:HB3	1.94	0.50
1:A:97:TYR:CB	1:A:100:ARG:HG2	2.42	0.50
1:A:557:LEU:C	1:A:578:ARG:HH21	2.20	0.50
1:A:364:ARG:HB3	1:A:366:ASP:OD1	2.12	0.50
1:A:523:ASP:OD1	1:A:526:LEU:HB2	2.12	0.49
1:A:195:ASP:OD1	1:A:197:ARG:HB3	2.13	0.49
1:A:495:GLU:HA	1:A:544:LEU:HD12	1.94	0.48
1:A:16:ALA:O	1:A:18:PRO:CD	2.60	0.48
1:A:590:ASP:OD1	1:A:591:GLN:HG3	2.13	0.48
1:A:321:ALA:HB2	1:A:344:CYS:SG	2.53	0.48
1:A:625:ILE:HD13	1:A:635:LEU:CD1	2.43	0.48
1:A:330:LEU:CB	1:A:331:PRO:HD3	2.43	0.48
1:A:507:ASP:OD1	1:A:509:HIS:HB2	2.14	0.48
1:A:86:PHE:HA	1:A:91:MET:HE2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:GLN:O	1:A:78:ASP:HB2	2.12	0.48
1:A:97:TYR:HB2	1:A:100:ARG:CG	2.44	0.48
1:A:363:GLN:HA	1:A:588:PHE:O	2.14	0.48
1:A:225:THR:CB	1:A:436:GLN:HE22	2.26	0.47
1:A:101:PHE:HE1	1:A:522:MET:HE2	1.80	0.47
1:A:101:PHE:CE1	1:A:522:MET:HE2	2.49	0.47
1:A:229:ASN:HA	1:A:242:THR:O	2.15	0.47
1:A:289:ILE:O	1:A:289:ILE:HD12	2.15	0.47
1:A:20:ASP:OD2	1:A:22:GLN:HB3	2.13	0.47
1:A:338:VAL:HG12	3:A:859:HOH:O	2.14	0.47
1:A:498:MET:CE	1:A:523:ASP:HB2	2.39	0.47
1:A:7:ASP:OD1	1:A:9:PRO:HD2	2.15	0.47
1:A:364:ARG:NH2	1:A:410:PRO:HG2	2.30	0.47
1:A:429:TYR:O	1:A:454:GLY:HA2	2.15	0.47
1:A:385:LEU:C	1:A:385:LEU:HD13	2.40	0.46
1:A:178:LEU:HA	1:A:262:MET:CE	2.45	0.46
1:A:238:GLN:HE21	1:A:238:GLN:HB3	1.59	0.46
1:A:469:VAL:O	1:A:472:ASP:HB2	2.15	0.45
1:A:539:THR:O	1:A:543:ARG:HG2	2.16	0.45
1:A:204:HIS:CD2	1:A:238:GLN:NE2	2.85	0.45
1:A:543:ARG:HH11	1:A:543:ARG:HB3	1.80	0.45
1:A:593:LEU:C	1:A:593:LEU:HD12	2.42	0.45
1:A:306:LEU:O	1:A:310:THR:HG23	2.17	0.45
1:A:597:LEU:CD1	1:A:623:GLY:H	2.29	0.45
1:A:6:ILE:HD13	1:A:6:ILE:N	2.31	0.45
1:A:318:VAL:HG13	1:A:345:HIS:CG	2.52	0.45
1:A:100:ARG:HB3	1:A:521:ALA:HA	1.98	0.45
1:A:204:HIS:CD2	1:A:238:GLN:HE22	2.35	0.45
1:A:510:ARG:HB3	1:A:516:TRP:CD1	2.52	0.45
1:A:291:THR:C	1:A:293:CYS:H	2.25	0.44
1:A:604:GLY:O	1:A:621:GLY:HA3	2.18	0.44
1:A:126:LEU:O	1:A:127:PRO:C	2.60	0.44
1:A:590:ASP:HA	1:A:629:PRO:HB3	2.00	0.44
1:A:20:ASP:HA	1:A:68:ARG:NH2	2.34	0.43
1:A:78:ASP:CG	1:A:315:PRO:HD2	2.43	0.43
1:A:240:MET:HE2	1:A:249:TRP:CE3	2.52	0.43
1:A:318:VAL:HG13	1:A:345:HIS:CB	2.47	0.43
1:A:88:GLN:HB2	1:A:90:HIS:CE1	2.54	0.43
1:A:132:ARG:HG3	1:A:184:ASP:OD1	2.17	0.43
1:A:269:LEU:CD2	3:A:819:HOH:O	2.65	0.43
1:A:214:TYR:CZ	1:A:289:ILE:HG23	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:HIS:HD2	1:A:238:GLN:NE2	2.17	0.43
1:A:605:TRP:O	1:A:635:LEU:HB3	2.18	0.43
1:A:124:HIS:HA	1:A:173:CYS:HB3	2.01	0.43
1:A:510:ARG:HD2	1:A:516:TRP:CD2	2.54	0.43
1:A:125:LEU:HD23	1:A:125:LEU:HA	1.88	0.43
1:A:534:LYS:HG3	1:A:534:LYS:O	2.19	0.42
1:A:164:ARG:O	1:A:167:GLU:HG2	2.19	0.42
1:A:314:ALA:O	1:A:317:VAL:HB	2.18	0.42
1:A:581:SER:O	1:A:636:GLN:HA	2.20	0.42
1:A:383:ALA:HB1	1:A:486:GLY:CA	2.50	0.42
1:A:602:VAL:HG23	1:A:603:ASP:N	2.34	0.42
1:A:522:MET:HG2	1:A:524:TRP:CE2	2.55	0.42
1:A:602:VAL:O	1:A:603:ASP:CB	2.65	0.42
1:A:351:THR:HG23	1:A:399:LEU:CD2	2.49	0.42
1:A:523:ASP:CG	1:A:526:LEU:HB2	2.43	0.42
1:A:587:ASN:C	1:A:587:ASN:HD22	2.27	0.42
1:A:81:ARG:HD3	1:A:85:TRP:CD1	2.55	0.42
1:A:205:PHE:N	1:A:238:GLN:HE21	2.18	0.42
1:A:355:ALA:CB	1:A:371:VAL:HG21	2.50	0.42
1:A:606:THR:CG2	1:A:620:ARG:HD2	2.49	0.42
1:A:31:TYR:O	1:A:35:ALA:HB2	2.20	0.42
1:A:158:LEU:C	1:A:158:LEU:CD2	2.93	0.42
1:A:112:VAL:N	1:A:113:PRO:CD	2.83	0.41
1:A:403:ALA:HA	1:A:412:PHE:CD1	2.55	0.41
2:B:1:GLC:O5	2:B:2:FRU:H11	2.19	0.41
1:A:132:ARG:HB2	1:A:182:ALA:HB1	2.01	0.41
1:A:327:MET:HE3	1:A:331:PRO:HD3	2.03	0.41
1:A:7:ASP:O	1:A:10:ALA:HB3	2.19	0.41
1:A:495:GLU:HA	1:A:544:LEU:CD1	2.50	0.41
1:A:590:ASP:OD1	1:A:590:ASP:C	2.63	0.41
1:A:318:VAL:HG13	1:A:345:HIS:CD2	2.55	0.41
1:A:443:HIS:HE1	3:A:932:HOH:O	2.04	0.41
1:A:527:ALA:O	1:A:530:ARG:HB3	2.21	0.41
1:A:331:PRO:N	1:A:332:PRO:HD2	2.36	0.41
1:A:574:PHE:O	1:A:586:HIS:HD2	2.04	0.41
1:A:429:TYR:CE2	1:A:473:ARG:CZ	3.04	0.41
1:A:529:GLN:C	1:A:531:HIS:N	2.78	0.40
1:A:222:PHE:N	1:A:223:PRO:HD3	2.37	0.40
1:A:463:ASP:OD1	1:A:463:ASP:C	2.64	0.40
1:A:237:ALA:O	1:A:238:GLN:HG2	2.21	0.40
1:A:543:ARG:HB3	1:A:543:ARG:NH1	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:582:PHE:CD1	1:A:582:PHE:C	2.99	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1024:HOH:O	3:A:1024:HOH:O[2_565]	1.68	0.52

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	612/644 (95%)	569 (93%)	35 (6%)	8 (1%)	9 8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	603	ASP
1	A	323	ALA
1	A	532	ASP
1	A	637	ARG
1	A	127	PRO
1	A	209	THR
1	A	227	PRO
1	A	511	GLN

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	474/491 (96%)	442 (93%)	32 (7%)	14	17

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ILE
1	A	20	ASP
1	A	77	LEU
1	A	118	LEU
1	A	136	ASN
1	A	161	LEU
1	A	164	ARG
1	A	172	LEU
1	A	178	LEU
1	A	199	LEU
1	A	207	ASP
1	A	208	ARG
1	A	251	LEU
1	A	264	LEU
1	A	269	LEU
1	A	274	VAL
1	A	289	ILE
1	A	317	VAL
1	A	338	VAL
1	A	350	SER
1	A	385	LEU
1	A	400	GLN
1	A	489	LEU
1	A	530	ARG
1	A	543	ARG
1	A	551	ARG
1	A	560	ASP
1	A	568	LEU
1	A	580	ASP
1	A	607	LEU
1	A	608	LEU
1	A	633	ARG

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	69	GLN
1	A	72	GLN
1	A	75	GLN
1	A	106	GLN
1	A	136	ASN
1	A	203	HIS
1	A	204	HIS
1	A	220	GLN
1	A	238	GLN
1	A	295	ASN
1	A	296	GLN
1	A	322	GLN
1	A	345	HIS
1	A	374	HIS
1	A	397	ASN
1	A	456	GLN
1	A	459	GLN
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 ⓘ

2 monosaccharides 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	GLC	B	1	2	11,11,12	1.12	1 (9%)	15,15,17	1.17	3 (20%)
2	FRU	B	2	2	11,12,12	2.82	1 (9%)	10,18,18	1.30	1 (10%)

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	B	1	2	-	0/2/19/22	0/1/1/1
2	FRU	B	2	2	-	5/5/24/24	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	FRU	O2-C2	-9.00	1.25	1.40
2	B	1	GLC	C2-C3	2.36	1.56	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	FRU	O2-C2-O5	3.32	115.70	109.33
2	B	1	GLC	C1-C2-C3	-2.35	106.22	109.64
2	B	1	GLC	C6-C5-C4	-2.33	107.29	113.02
2	B	1	GLC	C2-C3-C4	-2.19	107.02	110.86

There are no chirality outliers.

All (5) torsion outliers are listed below:

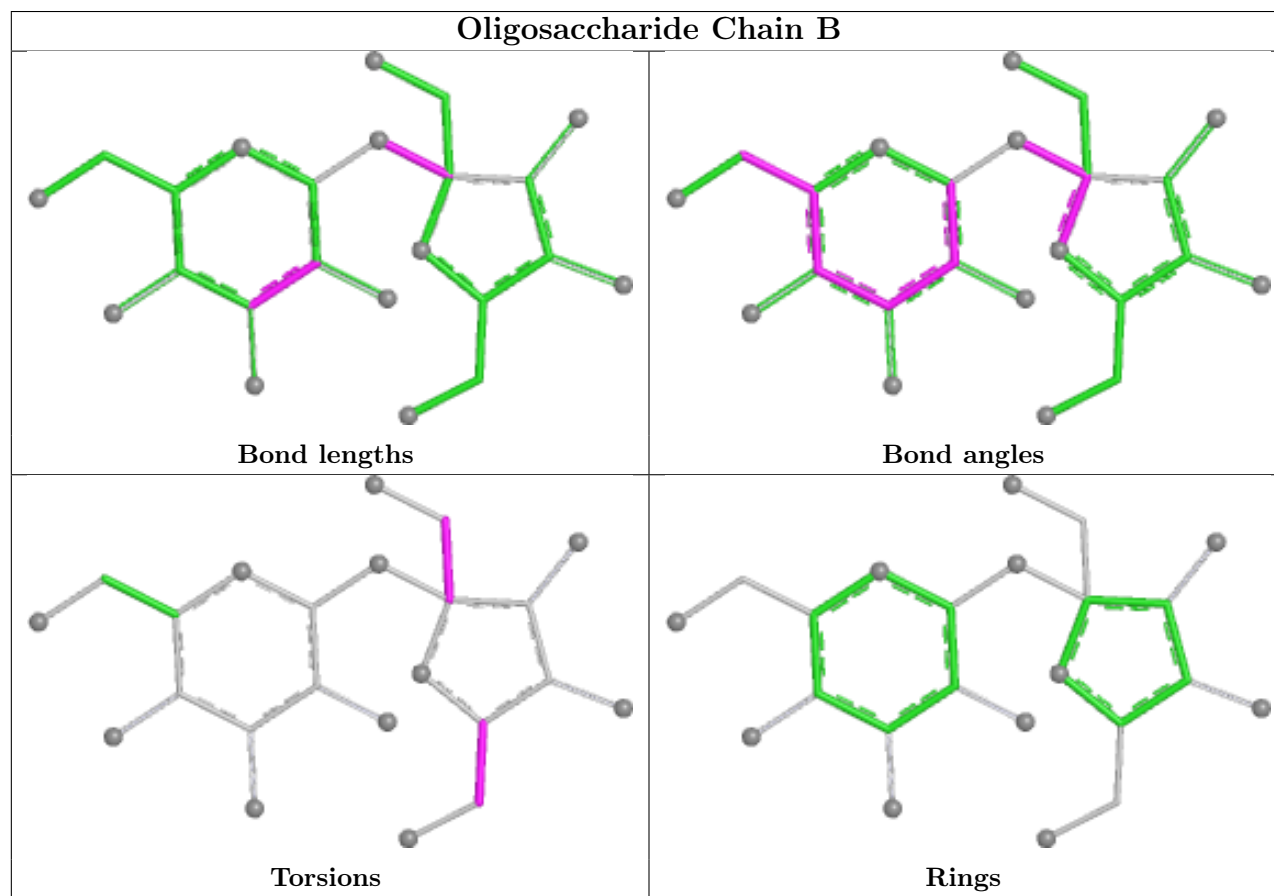
Mol	Chain	Res	Type	Atoms
2	B	2	FRU	O1-C1-C2-C3
2	B	2	FRU	O1-C1-C2-O2
2	B	2	FRU	O5-C5-C6-O6
2	B	2	FRU	C4-C5-C6-O6
2	B	2	FRU	O1-C1-C2-O5

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	FRU	1	0
2	B	1	GLC	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	618/644 (95%)	0.52	29 (4%) 36 33	20, 31, 52, 79	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	228	GLY	6.6
1	A	605	TRP	6.4
1	A	608	LEU	4.8
1	A	607	LEU	3.9
1	A	225	THR	3.3
1	A	621	GLY	3.2
1	A	638	GLY	3.2
1	A	224	HIS	3.2
1	A	68	ARG	3.1
1	A	606	THR	3.0
1	A	509	HIS	3.0
1	A	620	ARG	2.9
1	A	207	ASP	2.8
1	A	622	ASP	2.7
1	A	17	GLY	2.6
1	A	206	ALA	2.6
1	A	508	PRO	2.6
1	A	600	ILE	2.6
1	A	601	GLY	2.5
1	A	6	ILE	2.3
1	A	599	ALA	2.2
1	A	511	GLN	2.2
1	A	216	ALA	2.2
1	A	294	MET	2.1
1	A	227	PRO	2.1
1	A	213	ARG	2.1
1	A	443	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	426	PRO	2.1
1	A	210	VAL	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

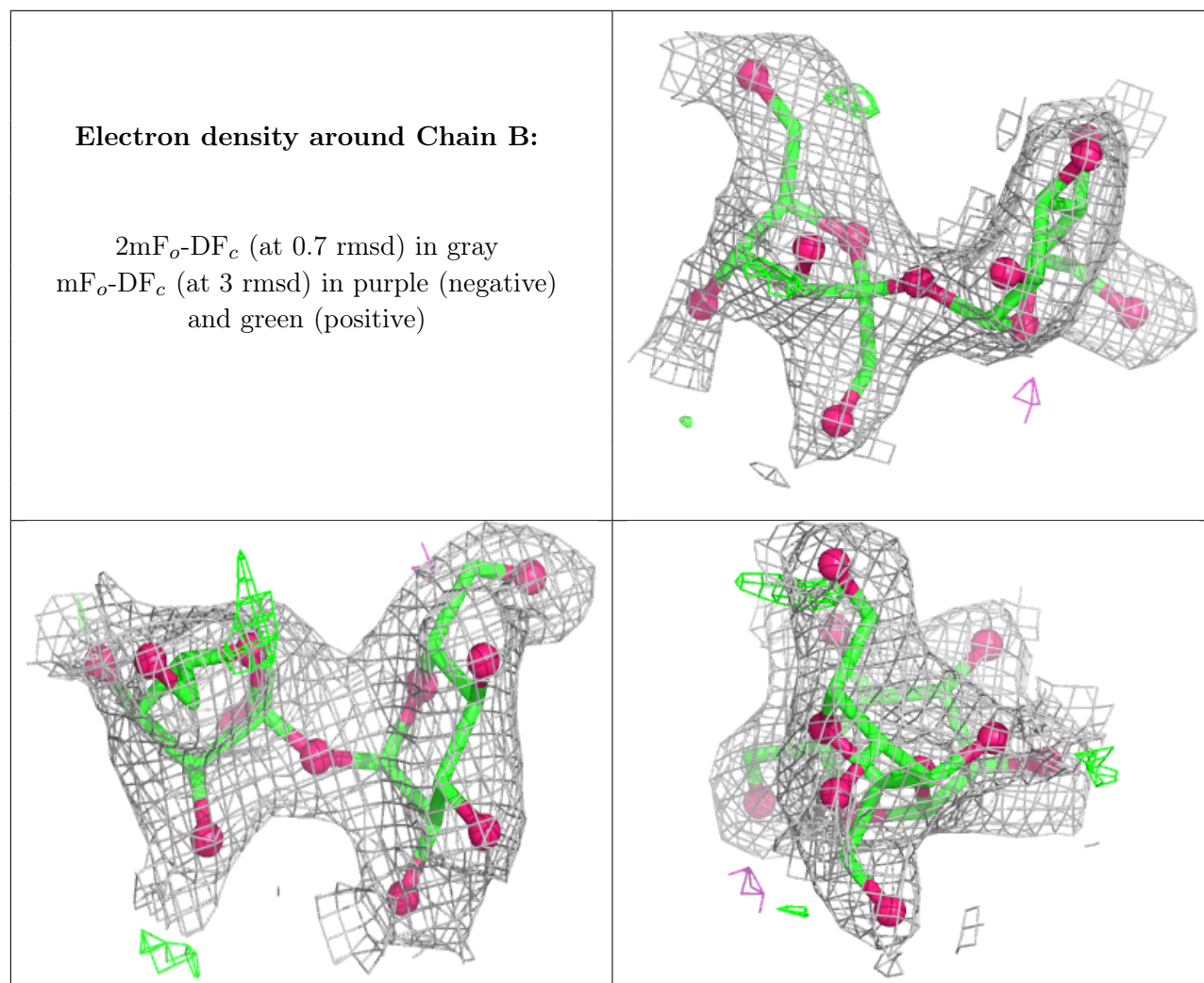
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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	FRU	B	2	12/12	0.77	0.15	45,50,52,53	0
2	GLC	B	1	11/12	0.85	0.11	39,40,43,43	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.