



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:21 PM UTC

PDB ID : 2HEZ / pdb_00002hez
Title : Bifidobacterium longum bile salt hydrolase
Authors : Suresh, C.G.; Kumar, R.S.; Brannigan, J.A.
Deposited on : 2006-06-22
Resolution : 2.50 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

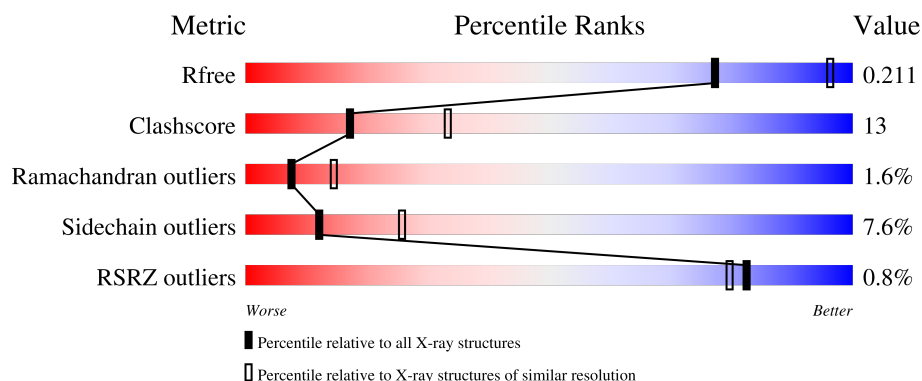
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	316	 73% 22% 5% .
1	B	316	 71% 22% 5% 2%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5164 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.

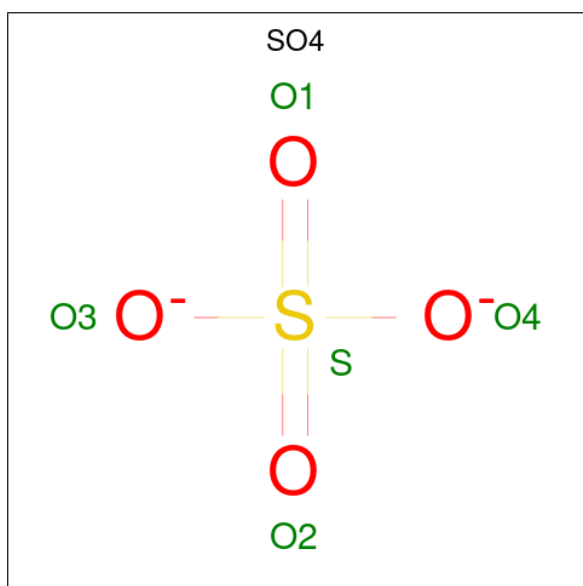
- Molecule 1 is a protein called Bile salt hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	14	0	0
			2466	1547	417	485	17			
1	B	316	Total	C	N	O	S	15	0	0
			2466	1547	417	485	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	OCS	CYS	modified residue	UNP Q9KK62
B	1	OCS	CYS	modified residue	UNP Q9KK62

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

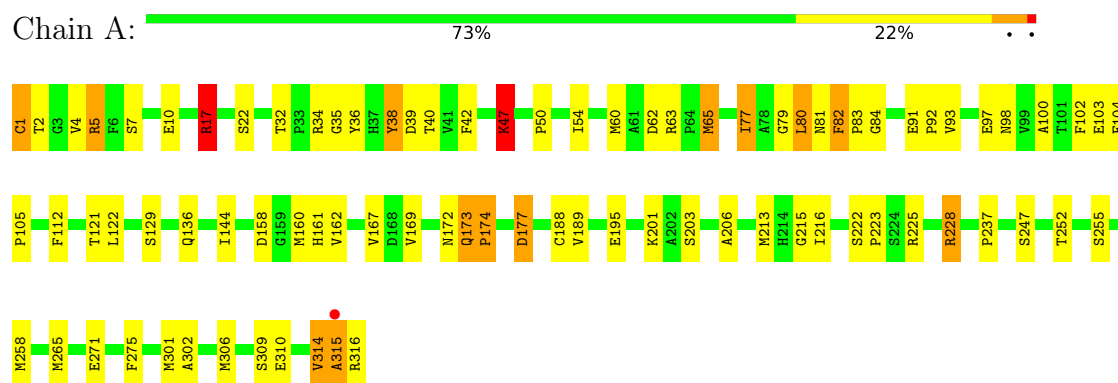
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	115	Total 115	O 115	0	0
3	B	112	Total 112	O 112	0	0

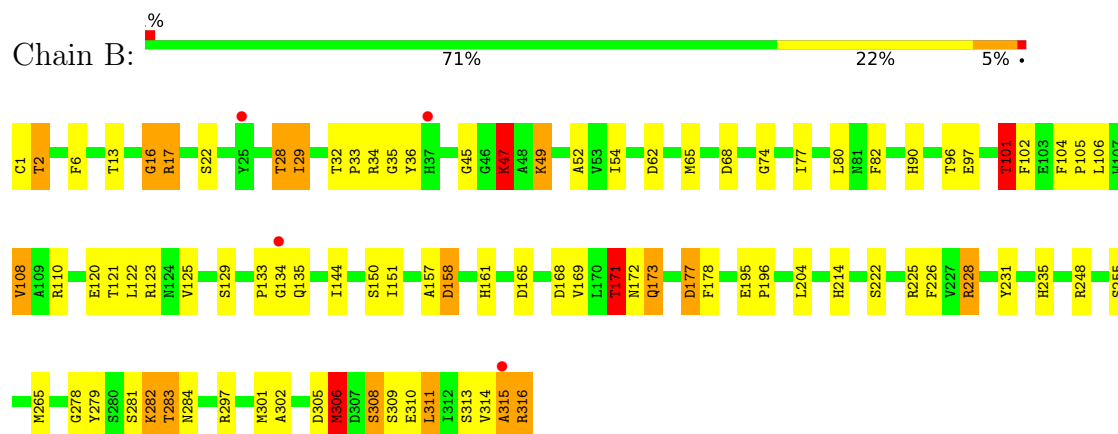
3 Residue-property plots

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: Bile salt hydrolase



- Molecule 1: Bile salt hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	125.24Å 125.24Å 117.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-2.50) 99.4 (20.00-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.174 , 0.216 0.173 , 0.211	Depositor DCC
R_{free} test set	1851 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	44.1	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5164	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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: OCS, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.40	5/2522 (0.2%)	1.34	21/3429 (0.6%)
1	B	1.43	8/2522 (0.3%)	1.39	20/3429 (0.6%)
All	All	1.41	13/5044 (0.3%)	1.37	41/6858 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	2	THR	CA-C	-6.96	1.38	1.52
1	B	17	ARG	CA-C	-6.92	1.44	1.52
1	A	215	GLY	C-O	-6.54	1.16	1.24
1	B	315	ALA	CA-C	-6.04	1.44	1.52
1	B	16	GLY	N-CA	6.00	1.51	1.45
1	B	157	ALA	CA-CB	5.83	1.62	1.53
1	B	177	ASP	CB-CG	5.66	1.66	1.52
1	B	82	PHE	CA-C	5.55	1.59	1.52
1	A	315	ALA	N-CA	5.47	1.53	1.46
1	A	206	ALA	CA-CB	-5.42	1.45	1.53
1	A	237	PRO	N-CA	5.23	1.53	1.47
1	B	110	ARG	C-O	-5.19	1.18	1.24
1	A	213	MET	SD-CE	-5.15	1.66	1.79

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2	THR	CA-CB-CG2	-11.39	91.14	110.50
1	B	2	THR	N-CA-C	-9.39	84.70	111.00
1	A	84	GLY	N-CA-C	-8.52	104.94	115.08
1	B	62	ASP	N-CA-C	-7.58	100.46	111.30
1	A	228	ARG	NE-CZ-NH2	-7.05	112.85	119.20
1	A	314	VAL	CA-C-O	-6.53	115.17	121.63
1	A	17	ARG	CG-CD-NE	-6.51	97.67	112.00
1	A	47	LYS	N-CA-C	-6.50	96.96	110.80
1	B	17	ARG	CG-CD-NE	-6.27	98.20	112.00
1	A	62	ASP	N-CA-C	-6.24	102.90	110.88
1	B	82	PHE	CA-C-N	-6.23	113.36	119.78
1	B	82	PHE	C-N-CA	-6.23	113.36	119.78
1	A	32	THR	CA-C-N	6.22	126.19	120.03
1	A	32	THR	C-N-CA	6.22	126.19	120.03
1	B	283	THR	CB-CA-C	-5.98	101.04	110.85
1	A	17	ARG	NE-CZ-NH2	-5.97	113.82	119.20
1	B	47	LYS	N-CA-C	-5.93	99.50	108.46
1	A	36	TYR	CA-C-N	-5.89	114.59	122.42
1	A	36	TYR	C-N-CA	-5.89	114.59	122.42
1	A	5	ARG	CG-CD-NE	-5.82	99.19	112.00
1	A	177	ASP	N-CA-CB	-5.81	101.58	110.12
1	A	38	TYR	CA-C-N	-5.56	114.80	122.30
1	A	38	TYR	C-N-CA	-5.56	114.80	122.30
1	A	167	VAL	N-CA-C	-5.49	107.63	112.90
1	B	282	LYS	N-CA-C	-5.45	104.74	111.33
1	B	171	THR	CB-CA-C	5.38	121.13	110.42
1	B	96	THR	N-CA-C	-5.31	101.91	110.14
1	B	158	ASP	N-CA-C	5.25	118.81	112.93
1	B	306	MET	N-CA-C	5.19	116.94	111.28
1	A	129	SER	CA-C-N	-5.16	115.60	122.72
1	A	129	SER	C-N-CA	-5.16	115.60	122.72
1	B	316	ARG	CD-NE-CZ	-5.14	117.21	124.40
1	B	228	ARG	NE-CZ-NH2	-5.13	114.58	119.20
1	A	247	SER	N-CA-C	-5.12	105.78	111.36
1	B	2	THR	N-CA-CB	-5.11	102.81	111.50
1	A	252	THR	N-CA-C	-5.11	105.79	111.36
1	B	151	ILE	N-CA-C	5.10	116.25	108.80
1	B	2	THR	CB-CA-C	5.10	120.32	109.10
1	B	195	GLU	CA-C-N	5.09	126.20	119.84
1	B	195	GLU	C-N-CA	5.09	126.20	119.84
1	A	177	ASP	CB-CA-C	-5.04	102.42	110.79

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	2	THR	Mainchain

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	2466	0	2288	48	0
1	B	2466	0	2288	71	1
2	B	5	0	0	0	0
3	A	115	0	0	2	1
3	B	112	0	0	5	0
All	All	5164	0	4576	119	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 13.

All (119) 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:49:LYS:CD	1:B:49:LYS:H	1.73	1.00
1:A:112:PHE:CE1	1:A:121:THR:HG21	2.00	0.96
1:B:173:GLN:NE2	1:B:173:GLN:H	1.67	0.93
1:B:49:LYS:H	1:B:49:LYS:HD3	1.32	0.91
1:A:172:ASN:HD22	1:A:225:ARG:HH22	1.22	0.87
1:B:172:ASN:HD22	1:B:225:ARG:HH22	1.21	0.87
1:A:310:GLU:HB2	3:A:425:HOH:O	1.74	0.87
1:B:101:THR:HG21	3:B:364:HOH:O	1.77	0.84
1:B:173:GLN:H	1:B:173:GLN:HE21	1.21	0.84
1:B:28:THR:HG23	1:B:29:ILE:O	1.78	0.83
1:B:235:HIS:HE1	3:B:349:HOH:O	1.61	0.81
1:A:22:SER:N	1:A:265:MET:HE1	1.96	0.80
1:A:112:PHE:CZ	1:A:121:THR:HG21	2.17	0.78
1:A:93:VAL:H	1:A:98:ASN:HD21	1.30	0.78
1:B:33:PRO:HA	1:B:306:MET:HE3	1.65	0.77
1:B:171:THR:HG23	1:B:172:ASN:H	1.48	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:ASN:ND2	1:B:225:ARG:HH22	1.82	0.77
1:B:65:MET:HE2	1:B:102:PHE:CD1	2.20	0.75
1:A:65:MET:HE1	1:A:102:PHE:CD1	2.23	0.73
1:A:172:ASN:ND2	1:A:225:ARG:HH22	1.84	0.73
1:B:171:THR:HG21	1:B:222:SER:O	1.89	0.72
1:B:35:GLY:O	1:B:309:SER:HB2	1.91	0.71
1:A:122:LEU:HD22	1:A:160:MET:HE1	1.73	0.70
1:B:101:THR:HG23	3:B:350:HOH:O	1.91	0.69
1:B:49:LYS:CD	1:B:49:LYS:N	2.52	0.69
1:B:158:ASP:OD1	1:B:161:HIS:HE1	1.76	0.68
1:B:33:PRO:HA	1:B:306:MET:CE	2.25	0.66
1:B:22:SER:HB3	1:B:265:MET:HE3	1.78	0.66
1:A:4:VAL:HG22	1:A:169:VAL:HG12	1.78	0.66
1:A:122:LEU:HB3	1:A:160:MET:HE1	1.77	0.66
1:A:173:GLN:H	1:A:173:GLN:NE2	1.95	0.65
1:B:28:THR:CG2	1:B:29:ILE:O	2.44	0.65
1:B:173:GLN:HE21	1:B:173:GLN:N	1.94	0.65
1:B:22:SER:CB	1:B:265:MET:HE3	2.26	0.65
1:A:93:VAL:H	1:A:98:ASN:ND2	1.95	0.64
1:B:32:THR:HG21	1:B:106:LEU:HD11	1.81	0.63
1:B:36:TYR:CE1	1:B:311:LEU:HD13	2.34	0.63
1:A:173:GLN:H	1:A:173:GLN:HE21	1.46	0.62
1:B:22:SER:N	1:B:265:MET:HE1	2.12	0.62
1:B:28:THR:HG21	1:B:313:SER:HB3	1.83	0.60
1:B:22:SER:N	1:B:265:MET:CE	2.65	0.59
1:A:172:ASN:HB3	1:A:173:GLN:NE2	2.18	0.58
1:B:49:LYS:H	1:B:49:LYS:HD2	1.63	0.58
1:B:45:GLY:N	1:B:97:GLU:OE1	2.33	0.58
1:B:34:ARG:NH1	1:B:306:MET:HB3	2.19	0.58
1:B:231:TYR:O	1:B:235:HIS:HD2	1.87	0.58
1:B:305:ASP:O	1:B:308:SER:HB3	2.04	0.57
1:B:77:ILE:HD12	1:B:144:ILE:HG12	1.86	0.55
1:B:74:GLY:HA2	1:B:279:TYR:OH	2.06	0.55
1:B:228:ARG:HD2	1:B:255:SER:O	2.07	0.55
1:A:225:ARG:NE	1:A:258:MET:HE2	2.21	0.55
1:A:228:ARG:HD2	1:A:255:SER:O	2.06	0.55
1:A:188:CYS:O	1:A:189:VAL:C	2.50	0.54
1:B:104:PHE:HB3	1:B:105:PRO:HD3	1.91	0.53
1:B:101:THR:HG22	3:B:371:HOH:O	2.10	0.51
1:B:311:LEU:O	1:B:311:LEU:HD23	2.10	0.51
1:A:112:PHE:CE1	1:A:121:THR:CG2	2.87	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:LYS:HG3	1:B:49:LYS:HD3	1.92	0.51
1:B:150:SER:HB2	1:B:168:ASP:OD1	2.11	0.51
1:B:283:THR:O	1:B:284:ASN:C	2.54	0.50
1:A:65:MET:HE1	1:A:102:PHE:CE1	2.45	0.50
1:B:120:GLU:OE2	1:B:120:GLU:HA	2.12	0.50
1:A:35:GLY:O	1:A:309:SER:HA	2.12	0.50
1:B:150:SER:HB3	1:B:165:ASP:HB3	1.94	0.49
1:A:79:GLY:C	1:A:80:LEU:HD23	2.38	0.49
1:B:120:GLU:O	1:B:123:ARG:HG3	2.13	0.49
1:B:54:ILE:HG22	1:B:301:MET:HE3	1.96	0.48
1:B:22:SER:H	1:B:265:MET:HE1	1.78	0.48
1:A:42:PHE:CD1	1:A:92:PRO:HG3	2.48	0.48
1:A:160:MET:HE2	1:A:162:VAL:CG2	2.43	0.48
1:B:178:PHE:CD2	1:B:178:PHE:C	2.91	0.48
1:B:101:THR:HG22	1:B:102:PHE:N	2.29	0.47
1:A:65:MET:CE	1:A:102:PHE:CD1	2.96	0.47
1:A:275:PHE:C	1:A:275:PHE:CD1	2.93	0.47
1:A:189:VAL:HG11	1:A:216:ILE:HG12	1.96	0.47
1:A:203:SER:HB3	3:A:360:HOH:O	2.14	0.47
1:A:100:ALA:O	1:A:103:GLU:HG2	2.15	0.47
1:A:158:ASP:OD1	1:A:161:HIS:HE1	1.96	0.47
1:B:311:LEU:HD22	3:B:424:HOH:O	2.15	0.47
1:A:77:ILE:HD12	1:A:144:ILE:HG12	1.96	0.46
1:B:6:PHE:CE1	1:B:248:ARG:HD3	2.49	0.46
1:B:171:THR:HG23	1:B:172:ASN:N	2.24	0.46
1:B:90:HIS:HD2	1:B:129:SER:O	1.97	0.46
1:B:65:MET:HE2	1:B:102:PHE:CE1	2.50	0.46
1:B:171:THR:CG2	1:B:172:ASN:H	2.24	0.46
1:A:122:LEU:HB3	1:A:160:MET:CE	2.45	0.45
1:B:104:PHE:O	1:B:108:VAL:HG13	2.17	0.45
1:A:22:SER:H	1:A:265:MET:HE1	1.75	0.45
1:B:121:THR:O	1:B:122:LEU:C	2.60	0.45
1:B:13:THR:HB	1:B:281:SER:HB3	1.98	0.45
1:A:39:ASP:O	1:A:63:ARG:NH2	2.50	0.44
1:A:34:ARG:HG2	1:A:50:PRO:O	2.16	0.44
1:B:121:THR:HG22	1:B:125:VAL:HG23	2.00	0.44
1:A:1:OCS:HB3	1:A:80:LEU:HB3	1.98	0.44
1:A:104:PHE:HB3	1:A:105:PRO:HD3	1.99	0.44
1:B:49:LYS:HD3	1:B:49:LYS:N	2.16	0.44
1:A:173:GLN:HB2	1:A:174:PRO:HA	2.00	0.43
1:B:28:THR:HG21	1:B:313:SER:CB	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:196:PRO:HA	1:B:204:LEU:O	2.19	0.43
1:B:34:ARG:HH12	1:B:306:MET:HB3	1.84	0.43
1:A:216:ILE:HG23	1:A:216:ILE:HD12	1.70	0.43
1:A:310:GLU:OE2	1:A:310:GLU:HA	2.18	0.43
1:A:301:MET:HG2	1:A:306:MET:HE3	2.01	0.42
1:B:101:THR:CG2	1:B:102:PHE:N	2.82	0.42
1:B:33:PRO:HG2	1:B:310:GLU:O	2.20	0.42
1:B:52:ALA:HB1	1:B:306:MET:HE2	2.02	0.42
1:A:222:SER:HB3	1:A:223:PRO:HD3	2.00	0.42
1:B:171:THR:CG2	1:B:172:ASN:N	2.83	0.42
1:B:49:LYS:N	1:B:49:LYS:HD2	2.29	0.42
1:B:231:TYR:O	1:B:235:HIS:CD2	2.71	0.41
1:B:52:ALA:CB	1:B:306:MET:HG3	2.50	0.41
1:B:171:THR:HG21	1:B:226:PHE:HB2	2.01	0.41
1:A:82:PHE:N	1:A:83:PRO:CD	2.83	0.41
1:A:54:ILE:HG22	1:A:301:MET:HE3	2.02	0.41
1:A:265:MET:HE2	1:A:271:GLU:HB2	2.03	0.41
1:A:2:THR:O	1:A:17:ARG:HA	2.20	0.41
1:A:38:TYR:CE2	1:A:40:THR:HG22	2.55	0.41
1:A:265:MET:HE2	1:A:271:GLU:CB	2.50	0.41
1:B:16:GLY:HA2	1:B:278:GLY:HA2	2.03	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)
1:B:214:HIS:CD2	3:A:416:HOH:O[5_555]	2.18	0.02

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	314/316 (99%)	295 (94%)	14 (4%)	5 (2%)	7	14
1	B	314/316 (99%)	290 (92%)	19 (6%)	5 (2%)	7	14
All	All	628/632 (99%)	585 (93%)	33 (5%)	10 (2%)	7	14

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	308	SER
1	A	47	LYS
1	A	302	ALA
1	B	101	THR
1	B	302	ALA
1	B	315	ALA
1	B	134	GLY
1	A	315	ALA
1	A	82	PHE
1	A	174	PRO

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	262/262 (100%)	243 (93%)	19 (7%)	13	27
1	B	262/262 (100%)	241 (92%)	21 (8%)	11	24
All	All	524/524 (100%)	484 (92%)	40 (8%)	12	26

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ARG
1	A	7	SER
1	A	10	GLU
1	A	17	ARG
1	A	47	LYS
1	A	60	MET

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Mol	Chain	Res	Type
1	A	65	MET
1	A	77	ILE
1	A	80	LEU
1	A	81	ASN
1	A	91	GLU
1	A	97	GLU
1	A	136	GLN
1	A	173	GLN
1	A	177	ASP
1	A	195	GLU
1	A	201	LYS
1	A	314	VAL
1	A	316	ARG
1	B	17	ARG
1	B	28	THR
1	B	29	ILE
1	B	47	LYS
1	B	49	LYS
1	B	68	ASP
1	B	80	LEU
1	B	101	THR
1	B	108	VAL
1	B	133	PRO
1	B	135	GLN
1	B	169	VAL
1	B	171	THR
1	B	173	GLN
1	B	177	ASP
1	B	282	LYS
1	B	297	ARG
1	B	306	MET
1	B	311	LEU
1	B	314	VAL
1	B	316	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	98	ASN
1	A	130	GLN
1	A	136	GLN

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Mol	Chain	Res	Type
1	A	161	HIS
1	A	172	ASN
1	A	173	GLN
1	A	185	ASN
1	A	233	ASN
1	A	245	ASN
1	A	269	GLN
1	A	284	ASN
1	B	18	ASN
1	B	90	HIS
1	B	111	ASN
1	B	161	HIS
1	B	172	ASN
1	B	173	GLN
1	B	235	HIS
1	B	245	ASN
1	B	257	GLN
1	B	284	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	OCS	B	1	1	6,8,9	2.60	1 (16%)	7,11,13	3.69	4 (57%)
1	OCS	A	1	1	6,8,9	2.12	2 (33%)	7,11,13	2.37	2 (28%)

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
1	OCS	B	1	1	-	2/4/7/9	-
1	OCS	A	1	1	-	1/4/7/9	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	OCS	CB-CA	-5.53	1.49	1.53
1	A	1	OCS	CB-CA	-3.76	1.50	1.53
1	A	1	OCS	OD3-SG	2.87	1.53	1.45

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	OCS	OD3-SG-CB	8.24	119.06	106.76
1	A	1	OCS	OD3-SG-CB	4.79	113.91	106.76
1	B	1	OCS	OD2-SG-OD3	-3.58	102.43	111.40
1	A	1	OCS	CA-CB-SG	-3.31	108.68	113.61
1	B	1	OCS	OD2-SG-OD1	2.53	117.74	111.40
1	B	1	OCS	OD1-SG-CB	-2.41	103.15	106.76

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	OCS	N-CA-CB-SG
1	B	1	OCS	N-CA-CB-SG
1	B	1	OCS	CA-CB-SG-OD1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1	OCS	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	SO4	B	317	-	4,4,4	0.45	0	6,6,6	1.38	1 (16%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	317	SO4	O3-S-O2	2.66	123.47	109.56

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	314/316 (99%)	-0.63	1 (0%) 90 87	23, 40, 64, 74	1 (0%)
1	B	314/316 (99%)	-0.52	4 (1%) 75 71	25, 42, 67, 77	2 (0%)
All	All	628/632 (99%)	-0.57	5 (0%) 82 80	23, 41, 66, 77	3 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	315	ALA	8.3
1	B	315	ALA	8.2
1	B	134	GLY	2.7
1	B	37	HIS	2.4
1	B	25	TYR	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	OCS	A	1	9/10	0.96	0.08	34,36,48,52	0
1	OCS	B	1	9/10	0.96	0.07	40,43,49,49	0

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	SO4	B	317	5/5	0.93	0.10	53,59,63,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.