



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:50 AM GMT

PDB ID : 2BXH
Title : HUMAN SERUM ALBUMIN COMPLEXED WITH INDOXYL SULFATE
Authors : Ghuman, J.; Zunszain, P.A.; Petitpas, I.; Bhattacharya, A.A.; Curry, S.
Deposited on : 2005-07-26
Resolution : 2.25 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

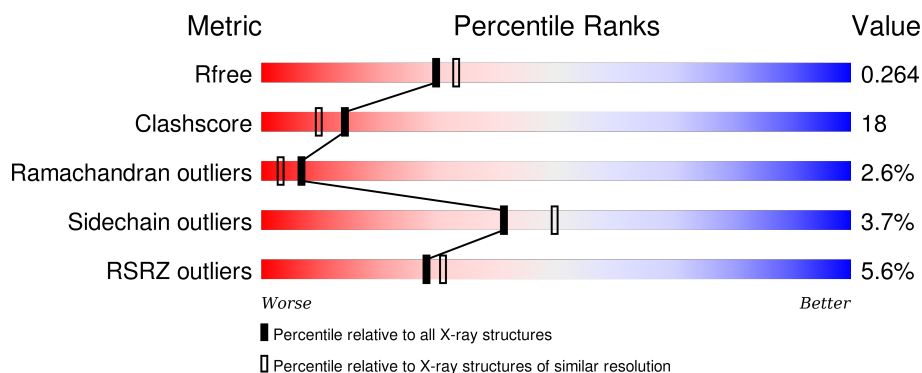
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	91344	1640 (2.28-2.24)
Clashscore	102246	1095 (2.26-2.26)
Ramachandran outliers	100387	1063 (2.26-2.26)
Sidechain outliers	100360	1063 (2.26-2.26)
RSRZ outliers	91569	1647 (2.28-2.24)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	585	<div> <div>5%</div> <div> <div></div> <div>64%</div> <div>30%</div> <div>• 5%</div> </div> </div>
1	B	585	<div> <div>5%</div> <div> <div></div> <div>64%</div> <div>29%</div> <div>• 5%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IOS	A	1002[A]	-	-	-	X
2	IOS	A	1002[B]	-	-	X	X
2	IOS	B	1002[A]	-	-	-	X
2	IOS	B	1002[B]	-	-	X	X

2 Entry composition [i](#)

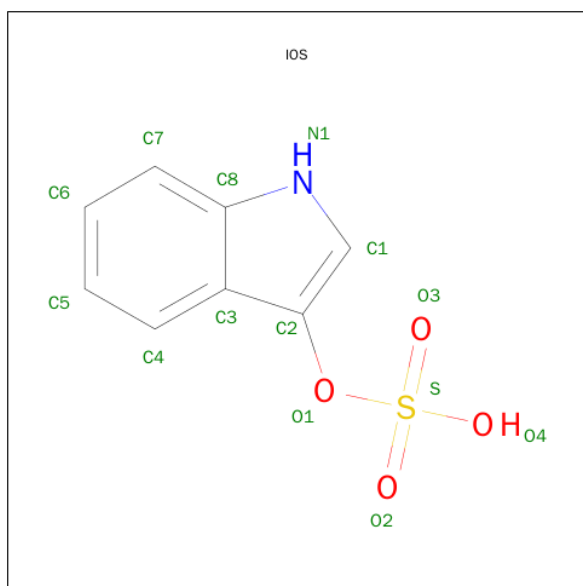
There are 3 unique types of molecules in this entry. The entry contains 8624 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4262	2699	714	809	40			
1	B	557	Total	C	N	O	S	0	0	0
			4224	2670	702	812	40			

- Molecule 2 is 3-SULFOOXY-1H-INDOLE (three-letter code: IOS) (formula: $C_8H_7NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			14	8	1	4	1		
2	A	1	Total	C	N	O	S	0	1
			28	16	2	8	2		
2	B	1	Total	C	N	O	S	0	0
			14	8	1	4	1		
2	B	1	Total	C	N	O	S	0	1
			28	16	2	8	2		

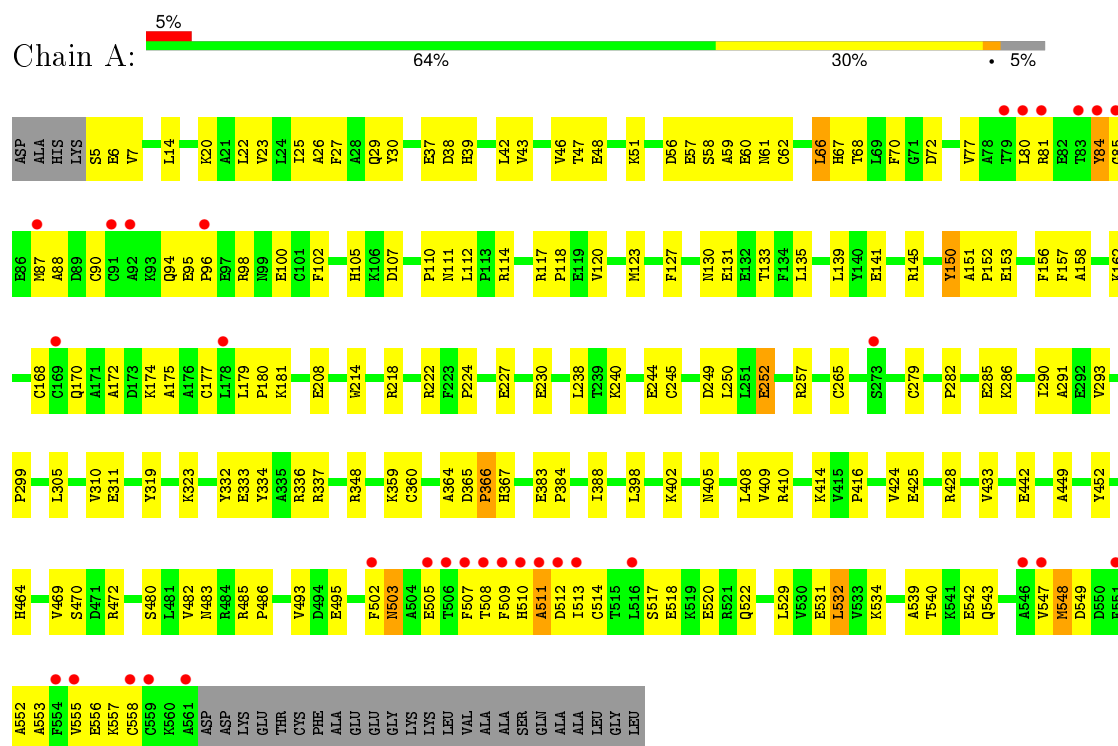
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	26	Total 26	O 26	0	0
3	B	28	Total 28	O 28	0	0

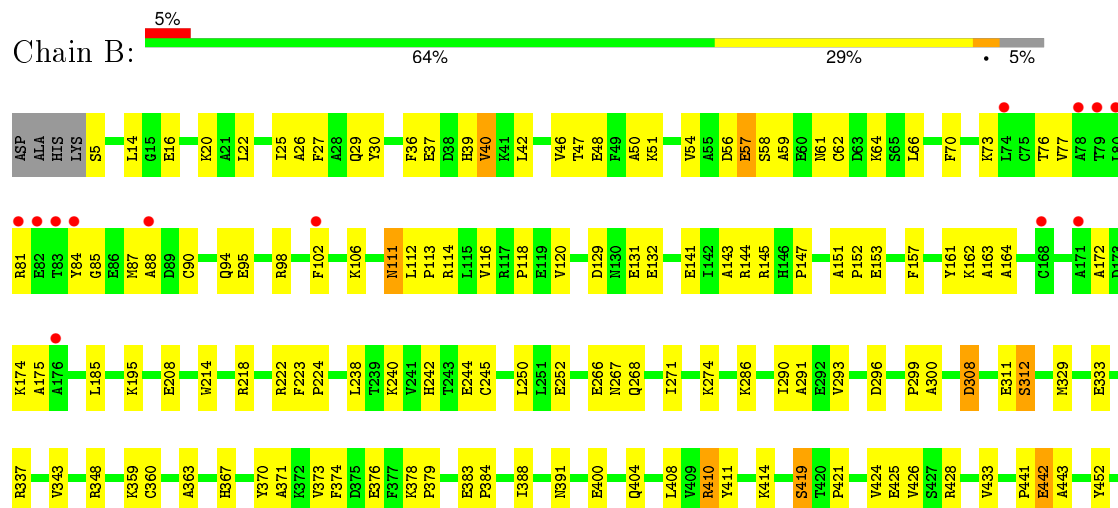
3 Residue-property plots

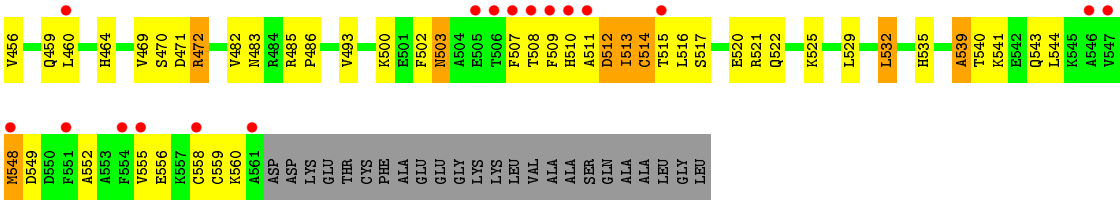
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: SERUM ALBUMIN



• Molecule 1: SERUM ALBUMIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.71Å 55.20Å 119.88Å 81.27° 91.09° 65.29°	Depositor
Resolution (Å)	35.03 – 2.25 36.31 – 2.25	Depositor EDS
% Data completeness (in resolution range)	98.6 (35.03-2.25) 92.6 (36.31-2.25)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.24Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.266 0.224 , 0.264	Depositor DCC
R_{free} test set	2973 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 53.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 59523 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8624	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.93% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: IOS

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	0.47	0/4347	0.64	0/5899
1	B	0.45	0/4308	0.63	2/5860 (0.0%)
All	All	0.46	0/8655	0.63	2/11759 (0.0%)

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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	472	ARG	NE-CZ-NH2	-8.39	116.11	120.30
1	B	472	ARG	NE-CZ-NH1	5.51	123.05	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	95	GLU	Mainchain

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	4262	0	4004	150	1
1	B	4224	0	3913	150	1
2	A	42	0	18	12	0
2	B	42	0	18	9	0
3	A	26	0	0	4	0
3	B	28	0	0	2	0
All	All	8624	0	7953	300	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 18.

All (300) 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:291:ALA:HB1	2:A:1002[B]:IOS:H4	1.31	1.11
1:B:348:ARG:HD2	1:B:483:ASN:ND2	1.85	0.91
1:B:540:THR:HB	1:B:543:GLN:HG2	1.53	0.91
1:A:214:TRP:HH2	2:A:1002[B]:IOS:H7	1.37	0.90
1:B:291:ALA:HB1	2:B:1002[B]:IOS:H4	1.56	0.88
1:B:517:SER:HB3	1:B:520:GLU:HG3	1.56	0.87
1:A:238:LEU:HD11	2:A:1002[B]:IOS:HA	1.56	0.86
1:B:410:ARG:HD3	3:B:2015:HOH:O	1.78	0.83
1:B:120:VAL:HG11	1:B:174:LYS:HB3	1.61	0.82
1:A:87:MET:HE1	1:A:105:HIS:HB3	1.61	0.82
1:A:265:CYS:SG	1:A:286:LYS:HE2	2.22	0.80
1:A:66:LEU:H	1:A:66:LEU:HD22	1.48	0.77
1:B:424:VAL:O	1:B:428:ARG:HG3	1.84	0.77
1:B:66:LEU:H	1:B:66:LEU:HD22	1.51	0.74
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.69	0.74
1:A:42:LEU:O	1:A:46:VAL:HG23	1.87	0.74
1:B:348:ARG:HD2	1:B:483:ASN:HD22	1.53	0.73
1:A:120:VAL:HG21	1:A:175:ALA:HA	1.70	0.73
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.69	0.72
1:A:87:MET:CE	1:A:105:HIS:HB3	2.21	0.70
1:B:510:HIS:O	1:B:513:ILE:HG13	1.92	0.70
1:A:424:VAL:O	1:A:428:ARG:HG3	1.91	0.69
1:A:240:LYS:O	1:A:244:GLU:HG3	1.92	0.69
1:A:333:GLU:HA	1:A:336:ARG:HD2	1.75	0.69
1:A:218:ARG:NH2	1:A:222:ARG:HH21	1.90	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:GLU:O	1:A:157:PHE:HD1	1.77	0.68
1:A:442:GLU:CD	1:A:442:GLU:H	1.97	0.68
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.76	0.68
1:A:179:LEU:HB2	1:A:180:PRO:HD3	1.75	0.68
1:B:14:LEU:HD13	1:B:22:LEU:HD12	1.76	0.67
1:B:151:ALA:HB3	1:B:152:PRO:HD3	1.77	0.67
1:B:61:ASN:HB3	1:B:64:LYS:HD3	1.76	0.66
1:A:348:ARG:HG3	1:A:482:VAL:HG12	1.78	0.66
1:A:156:PHE:HE1	1:A:285:GLU:HG3	1.61	0.66
1:B:50:ALA:O	1:B:54:VAL:HG23	1.96	0.65
1:B:367:HIS:HA	1:B:370:TYR:CZ	2.31	0.65
1:A:365:ASP:N	1:A:366:PRO:HD3	2.11	0.65
1:B:286:LYS:O	1:B:290:ILE:HG13	1.97	0.65
1:A:348:ARG:HG3	1:A:482:VAL:CG1	2.27	0.64
1:A:305:LEU:HD21	1:A:333:GLU:HB3	1.79	0.64
1:B:552:ALA:O	1:B:555:VAL:HG12	1.98	0.63
1:A:398:LEU:HB3	1:A:402:LYS:HB3	1.79	0.63
1:B:391:ASN:OD1	1:B:410:ARG:NH2	2.30	0.63
1:A:503:ASN:HD21	1:A:505:GLU:HB2	1.63	0.63
1:B:274:LYS:HE3	1:B:296:ASP:HA	1.80	0.63
1:B:483:ASN:C	1:B:486:PRO:HD2	2.18	0.63
1:A:139:LEU:HD21	1:A:158:ALA:HB2	1.80	0.62
1:A:87:MET:O	1:A:90:CYS:HB2	1.98	0.62
1:B:384:PRO:O	1:B:388:ILE:HG12	1.99	0.62
1:B:348:ARG:HD3	1:B:482:VAL:HG12	1.81	0.61
1:B:539:ALA:HA	1:B:543:GLN:HE21	1.65	0.61
1:A:168:CYS:SG	1:A:177:CYS:C	2.78	0.61
1:B:141:GLU:HA	1:B:141:GLU:OE1	2.00	0.61
1:B:224:PRO:HB2	1:B:299:PRO:HD3	1.83	0.61
1:B:94:GLN:O	1:B:98:ARG:N	2.31	0.61
1:B:290:ILE:O	1:B:293:VAL:HG12	2.01	0.61
1:A:384:PRO:O	1:A:388:ILE:HG12	2.00	0.60
1:B:378:LYS:CB	1:B:379:PRO:HD3	2.30	0.60
1:A:414:LYS:O	1:A:472:ARG:NH1	2.35	0.60
1:B:464:HIS:HE1	1:B:470:SER:H	1.49	0.60
1:A:60:GLU:HG3	1:A:61:ASN:OD1	2.02	0.60
1:A:66:LEU:O	1:A:70:PHE:HD1	1.85	0.60
1:A:540:THR:HG23	1:A:542:GLU:H	1.67	0.60
1:A:26:ALA:O	1:A:30:TYR:HD2	1.85	0.59
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.84	0.59
1:B:132:GLU:OE1	1:B:132:GLU:N	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:ARG:CZ	1:B:88:ALA:HB3	2.32	0.59
1:A:30:TYR:CD1	1:A:102:PHE:HB3	2.38	0.59
1:B:539:ALA:HA	1:B:543:GLN:NE2	2.18	0.59
1:B:333:GLU:OE1	1:B:337:ARG:NH2	2.36	0.59
1:A:410:ARG:HD2	3:A:2019:HOH:O	2.03	0.58
1:B:483:ASN:O	1:B:486:PRO:HD2	2.04	0.58
1:A:224:PRO:HB2	1:A:299:PRO:HD3	1.86	0.58
1:A:513:ILE:HD12	1:A:514:CYS:SG	2.44	0.58
1:B:131:GLU:HB3	1:B:132:GLU:OE1	2.03	0.58
1:A:81:ARG:HG2	1:A:88:ALA:CB	2.34	0.57
1:A:120:VAL:HG11	1:A:174:LYS:HB3	1.86	0.57
1:A:214:TRP:CH2	2:A:1002[B]:IOS:H7	2.29	0.57
1:A:532:LEU:HD11	1:A:547:VAL:CG1	2.35	0.57
1:A:290:ILE:O	1:A:293:VAL:HG12	2.04	0.57
1:A:508:THR:HG22	1:A:510:HIS:H	1.70	0.57
1:B:214:TRP:HH2	2:B:1002[A]:IOS:H7	1.69	0.57
1:B:224:PRO:HD2	1:B:296:ASP:HB3	1.87	0.57
1:B:76:THR:O	1:B:76:THR:HG22	2.05	0.56
1:B:16:GLU:O	1:B:20:LYS:HG3	2.05	0.56
1:A:26:ALA:HB2	1:A:250:LEU:HD12	1.88	0.56
1:B:36:PHE:O	1:B:40:VAL:HG23	2.04	0.56
1:B:464:HIS:CE1	1:B:470:SER:H	2.23	0.56
1:B:529:LEU:HD22	1:B:548:MET:HG3	1.88	0.55
1:A:95:GLU:OE1	1:A:96:PRO:HA	2.06	0.55
1:A:5:SER:HA	1:A:62:CYS:O	2.06	0.55
1:A:238:LEU:HD21	2:A:1002[B]:IOS:HA	1.89	0.55
1:B:145:ARG:HH11	1:B:145:ARG:HG3	1.71	0.55
1:B:400:GLU:O	1:B:404:GLN:HG3	2.07	0.55
1:B:56:ASP:C	1:B:58:SER:H	2.10	0.55
1:B:433:VAL:HG22	1:B:452:TYR:CD1	2.41	0.54
1:A:127:PHE:CE1	1:A:131:GLU:HG2	2.42	0.54
1:B:153:GLU:O	1:B:157:PHE:HD1	1.90	0.54
1:A:27:PHE:CE1	1:A:70:PHE:HD2	2.25	0.54
1:A:282:PRO:O	1:A:286:LYS:HB2	2.07	0.54
1:A:6:GLU:CG	1:A:66:LEU:HD21	2.38	0.54
1:A:364:ALA:C	1:A:366:PRO:HD3	2.28	0.54
1:B:266:GLU:C	1:B:268:GLN:H	2.11	0.54
1:A:30:TYR:HD1	1:A:102:PHE:HB3	1.73	0.54
1:A:333:GLU:OE1	1:A:337:ARG:NH2	2.41	0.54
1:A:553:ALA:O	1:A:556:GLU:HB3	2.08	0.54
1:A:464:HIS:HE1	1:A:470:SER:H	1.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:ASP:O	1:A:110:PRO:HD3	2.07	0.54
1:B:373:VAL:HG13	1:B:374:PHE:HD1	1.73	0.53
1:B:500:LYS:O	1:B:535:HIS:ND1	2.34	0.53
1:B:456:VAL:O	1:B:459:GLN:HB2	2.07	0.53
1:B:521:ARG:HD3	3:B:2027:HOH:O	2.08	0.53
1:A:130:ASN:ND2	1:A:133:THR:HB	2.23	0.53
1:B:522:GLN:HA	1:B:525:LYS:HE2	1.91	0.53
1:A:222:ARG:NH2	2:A:1002[A]:IOS:O3	2.42	0.53
1:A:84:TYR:HB2	1:A:87:MET:HB2	1.90	0.53
1:A:552:ALA:O	1:A:555:VAL:HG12	2.09	0.53
1:A:540:THR:H	1:A:543:GLN:CG	2.22	0.53
1:A:56:ASP:C	1:A:58:SER:H	2.12	0.52
1:A:25:ILE:O	1:A:29:GLN:HG3	2.09	0.52
1:A:7:VAL:HG23	3:A:1998:HOH:O	2.08	0.52
1:A:503:ASN:ND2	1:A:505:GLU:HB2	2.25	0.52
1:B:161:TYR:O	1:B:164:ALA:HB3	2.09	0.52
1:B:441:PRO:O	1:B:443:ALA:N	2.42	0.52
1:B:61:ASN:O	1:B:64:LYS:HB2	2.09	0.52
1:B:485:ARG:HB3	1:B:486:PRO:HD3	1.91	0.52
1:A:107:ASP:OD2	1:A:110:PRO:HA	2.09	0.52
1:A:483:ASN:C	1:A:486:PRO:HD2	2.30	0.51
1:A:291:ALA:O	2:A:1002[A]:IOS:H5	2.10	0.51
1:A:529:LEU:HD13	1:A:548:MET:CG	2.41	0.51
1:B:308:ASP:N	1:B:308:ASP:OD2	2.44	0.51
1:A:168:CYS:O	1:A:174:LYS:HG2	2.11	0.51
1:B:141:GLU:OE1	1:B:144:ARG:HD3	2.11	0.51
1:B:238:LEU:HD21	2:B:1002[B]:IOS:HA	1.93	0.51
1:B:120:VAL:HG21	1:B:175:ALA:HA	1.92	0.51
1:B:66:LEU:O	1:B:70:PHE:HD1	1.94	0.51
1:B:513:ILE:HD12	1:B:513:ILE:H	1.76	0.50
1:B:42:LEU:O	1:B:46:VAL:HG23	2.10	0.50
1:B:25:ILE:O	1:B:29:GLN:HG3	2.11	0.50
1:A:508:THR:HG22	1:A:509:PHE:N	2.27	0.50
1:B:511:ALA:O	1:B:513:ILE:HD12	2.11	0.50
1:A:81:ARG:NH1	1:A:88:ALA:HB3	2.27	0.50
1:A:48:GLU:O	1:A:51:LYS:HB2	2.11	0.50
1:A:252:GLU:OE1	1:A:252:GLU:N	2.30	0.50
1:B:508:THR:HG22	1:B:509:PHE:H	1.76	0.50
1:A:540:THR:HG22	1:A:543:GLN:HG2	1.93	0.49
1:A:56:ASP:O	1:A:58:SER:N	2.40	0.49
1:B:508:THR:HG22	1:B:509:PHE:N	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:GLU:HG2	1:A:66:LEU:HD21	1.93	0.49
1:B:329:MET:O	1:B:333:GLU:HG2	2.12	0.49
1:B:30:TYR:HD1	1:B:102:PHE:HB3	1.77	0.49
1:B:87:MET:O	1:B:90:CYS:HB2	2.12	0.49
1:A:39:HIS:O	1:A:43:VAL:HG23	2.12	0.49
1:A:388:ILE:HD12	1:A:449:ALA:CB	2.43	0.49
1:B:540:THR:O	1:B:541:LYS:C	2.50	0.49
1:B:222:ARG:O	1:B:224:PRO:HD3	2.13	0.48
1:B:30:TYR:CD1	1:B:102:PHE:HB3	2.47	0.48
1:A:94:GLN:O	1:A:98:ARG:N	2.44	0.48
1:B:512:ASP:C	1:B:515:THR:HG22	2.33	0.48
1:A:214:TRP:CH2	2:A:1002[A]:IOS:N1	2.81	0.48
1:B:120:VAL:HG21	1:B:175:ALA:CA	2.43	0.48
1:A:365:ASP:N	1:A:366:PRO:CD	2.74	0.48
1:A:177:CYS:O	1:A:181:LYS:HG3	2.14	0.48
1:B:311:GLU:O	1:B:312:SER:O	2.32	0.48
1:A:517:SER:HB3	1:A:520:GLU:HG3	1.96	0.48
1:B:5:SER:HA	1:B:62:CYS:O	2.14	0.48
1:A:319:TYR:CE1	1:A:323:LYS:HB2	2.49	0.48
1:A:238:LEU:HD21	2:A:1002[B]:IOS:C1	2.44	0.47
1:B:511:ALA:C	1:B:513:ILE:H	2.18	0.47
1:B:311:GLU:O	1:B:367:HIS:HE1	1.97	0.47
1:A:359:LYS:CG	1:A:360:CYS:N	2.77	0.47
1:B:556:GLU:O	1:B:559:CYS:HB2	2.14	0.47
1:A:531:GLU:HA	1:A:531:GLU:OE1	2.14	0.47
1:B:47:THR:O	1:B:50:ALA:HB3	2.14	0.47
1:B:111:ASN:O	1:B:112:LEU:HD23	2.15	0.47
1:A:23:VAL:HG22	1:A:70:PHE:HE2	1.79	0.47
1:A:408:LEU:HD21	1:A:424:VAL:HA	1.96	0.47
1:A:509:PHE:O	1:A:510:HIS:C	2.53	0.47
1:B:549:ASP:O	1:B:552:ALA:HB3	2.15	0.47
1:B:208:GLU:H	1:B:208:GLU:CD	2.16	0.47
1:A:214:TRP:HH2	2:A:1002[A]:IOS:N1	2.12	0.47
1:A:81:ARG:HG2	1:A:88:ALA:HB3	1.97	0.47
1:B:529:LEU:HD22	1:B:548:MET:CG	2.45	0.47
1:B:208:GLU:N	1:B:208:GLU:OE1	2.36	0.47
1:B:113:PRO:O	1:B:114:ARG:C	2.53	0.47
1:A:502:PHE:CD2	1:A:503:ASN:N	2.83	0.46
1:A:410:ARG:CD	3:A:2019:HOH:O	2.61	0.46
1:A:442:GLU:CD	1:A:442:GLU:N	2.65	0.46
1:B:441:PRO:O	1:B:442:GLU:C	2.53	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:529:LEU:HD13	1:A:548:MET:HG2	1.96	0.46
1:A:540:THR:H	1:A:543:GLN:HG2	1.79	0.46
1:A:405:ASN:O	1:A:409:VAL:HG23	2.14	0.46
1:A:66:LEU:CD2	1:A:66:LEU:H	2.25	0.46
1:A:549:ASP:O	1:A:552:ALA:N	2.49	0.46
1:B:66:LEU:N	1:B:66:LEU:HD22	2.25	0.46
1:B:367:HIS:O	1:B:371:ALA:HB2	2.16	0.46
1:B:37:GLU:CD	1:B:37:GLU:H	2.18	0.46
1:B:66:LEU:H	1:B:66:LEU:CD2	2.24	0.46
1:B:120:VAL:HG21	1:B:175:ALA:HB2	1.98	0.45
1:A:532:LEU:HD11	1:A:547:VAL:HG11	1.99	0.45
1:A:141:GLU:OE1	1:A:141:GLU:HA	2.15	0.45
1:A:511:ALA:O	1:A:513:ILE:N	2.38	0.45
1:B:419:SER:OG	1:B:421:PRO:HD2	2.17	0.45
1:B:540:THR:H	1:B:543:GLN:HG3	1.82	0.45
1:B:195:LYS:HD3	2:B:1002[B]:IOS:C6	2.47	0.45
1:A:152:PRO:CB	1:A:257:ARG:HH11	2.28	0.45
1:B:376:GLU:O	1:B:379:PRO:HD2	2.17	0.45
1:A:279:CYS:O	1:A:286:LYS:HE3	2.17	0.45
1:B:291:ALA:HB1	2:B:1002[A]:IOS:H5	1.97	0.45
1:A:111:ASN:O	1:A:112:LEU:HD23	2.17	0.45
1:A:416:PRO:O	1:A:534:LYS:HE2	2.17	0.45
1:A:279:CYS:HA	1:A:286:LYS:HE2	1.99	0.44
1:B:370:TYR:C	1:B:370:TYR:CD1	2.90	0.44
1:A:513:ILE:HD13	1:A:555:VAL:HG23	1.99	0.44
1:A:310:VAL:HG23	1:A:311:GLU:N	2.31	0.44
1:B:408:LEU:HD21	1:B:424:VAL:HA	1.98	0.44
1:B:56:ASP:O	1:B:58:SER:N	2.50	0.44
1:B:513:ILE:CD1	1:B:514:CYS:SG	3.06	0.44
1:B:222:ARG:C	1:B:224:PRO:HD3	2.38	0.44
1:B:464:HIS:CE1	1:B:469:VAL:H	2.35	0.44
1:A:20:LYS:HD2	1:A:47:THR:HG21	1.99	0.44
1:B:426:VAL:HG21	1:B:460:LEU:HB2	1.99	0.44
1:A:23:VAL:HG22	1:A:70:PHE:CE2	2.51	0.44
1:B:106:LYS:HD3	1:B:147:PRO:HB2	1.98	0.44
1:B:214:TRP:HH2	2:B:1002[B]:IOS:H7	1.83	0.44
1:B:66:LEU:O	1:B:70:PHE:CD1	2.71	0.44
1:B:162:LYS:C	1:B:164:ALA:H	2.20	0.44
1:A:6:GLU:HG3	1:A:66:LEU:HD21	2.00	0.44
1:A:14:LEU:HD13	1:A:22:LEU:HD12	1.99	0.44
1:A:433:VAL:HG22	1:A:452:TYR:CD1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:TRP:CD1	1:B:343:VAL:HG11	2.52	0.44
1:A:279:CYS:HA	1:A:286:LYS:CE	2.48	0.44
1:A:6:GLU:N	3:A:1998:HOH:O	2.51	0.44
1:B:521:ARG:O	1:B:525:LYS:HG3	2.18	0.44
1:B:540:THR:O	1:B:544:LEU:N	2.51	0.43
1:A:348:ARG:CG	1:A:482:VAL:HG12	2.47	0.43
1:B:88:ALA:C	1:B:90:CYS:N	2.71	0.43
1:A:117:ARG:HB2	1:A:123:MET:CE	2.48	0.43
1:B:116:VAL:O	1:B:116:VAL:HG13	2.17	0.43
1:A:6:GLU:HG2	1:A:66:LEU:CD2	2.48	0.43
1:A:135:LEU:HD11	1:A:162:LYS:HB2	2.00	0.43
1:B:507:PHE:CD1	1:B:507:PHE:O	2.71	0.43
1:A:238:LEU:CD1	2:A:1002[B]:IOS:HA	2.38	0.43
1:B:517:SER:O	1:B:520:GLU:N	2.51	0.43
1:B:512:ASP:CA	1:B:515:THR:HG22	2.49	0.43
1:A:503:ASN:ND2	1:A:505:GLU:H	2.17	0.42
1:A:120:VAL:HG11	1:A:174:LYS:CB	2.49	0.42
1:B:81:ARG:NH1	1:B:88:ALA:HB3	2.35	0.42
1:B:48:GLU:O	1:B:51:LYS:N	2.52	0.42
1:A:332:TYR:OH	1:A:336:ARG:NH2	2.53	0.42
1:B:26:ALA:HB2	1:B:250:LEU:HD12	2.02	0.42
1:B:27:PHE:HB3	1:B:39:HIS:CE1	2.54	0.42
1:A:150:TYR:HE2	2:A:1002[B]:IOS:O3	2.03	0.42
1:B:240:LYS:O	1:B:244:GLU:HG3	2.20	0.42
1:A:145:ARG:HG3	1:A:145:ARG:HH11	1.84	0.42
1:B:218:ARG:HD2	1:B:218:ARG:HA	1.91	0.42
1:A:513:ILE:C	1:A:513:ILE:HD12	2.39	0.42
1:A:507:PHE:O	1:A:507:PHE:CD1	2.73	0.42
1:A:557:LYS:O	1:A:558:CYS:C	2.58	0.42
1:A:130:ASN:ND2	1:A:133:THR:CB	2.83	0.41
1:B:48:GLU:O	1:B:51:LYS:HB2	2.20	0.41
1:A:518:GLU:OE2	1:A:522:GLN:HG3	2.19	0.41
1:B:29:GLN:HG2	1:B:143:ALA:O	2.19	0.41
1:A:100:GLU:HA	1:A:100:GLU:OE1	2.20	0.41
1:A:513:ILE:CD1	1:A:514:CYS:SG	3.08	0.41
1:B:291:ALA:HB1	2:B:1002[B]:IOS:C4	2.41	0.41
1:B:218:ARG:HD3	1:B:343:VAL:HG21	2.02	0.41
1:B:57:GLU:CD	1:B:57:GLU:H	2.23	0.41
1:B:333:GLU:O	1:B:337:ARG:NE	2.53	0.41
1:B:525:LYS:O	1:B:548:MET:HE1	2.21	0.41
1:A:311:GLU:HA	1:A:367:HIS:HE1	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:348:ARG:CD	1:B:483:ASN:ND2	2.72	0.41
1:B:88:ALA:O	1:B:90:CYS:N	2.54	0.41
1:A:508:THR:CG2	1:A:509:PHE:N	2.84	0.41
1:A:464:HIS:CE1	1:A:470:SER:H	2.37	0.41
1:A:359:LYS:CG	1:A:360:CYS:H	2.34	0.41
1:A:68:THR:O	1:A:72:ASP:OD1	2.38	0.41
1:B:120:VAL:HG21	1:B:175:ALA:CB	2.51	0.41
1:B:411:TYR:HA	1:B:414:LYS:HD3	2.03	0.41
1:B:222:ARG:HG2	1:B:223:PHE:CE1	2.56	0.40
1:B:56:ASP:C	1:B:58:SER:N	2.75	0.40
1:B:532:LEU:O	1:B:535:HIS:HB3	2.21	0.40
1:A:227:GLU:O	1:A:230:GLU:N	2.50	0.40
1:A:67:HIS:NE2	1:A:249:ASP:OD1	2.54	0.40
1:A:27:PHE:HE1	1:A:70:PHE:HD2	1.66	0.40
1:A:168:CYS:SG	1:A:177:CYS:O	2.80	0.40
1:B:145:ARG:HG3	1:B:145:ARG:NH1	2.35	0.40
1:B:512:ASP:O	1:B:515:THR:HG22	2.21	0.40
1:B:185:LEU:HD23	1:B:185:LEU:HA	1.91	0.40
1:B:516:LEU:HB3	1:B:520:GLU:HB2	2.03	0.40
1:B:238:LEU:HD11	2:B:1002[B]:IOS:O4	2.21	0.40
1:B:162:LYS:C	1:B:164:ALA:N	2.75	0.40
1:B:502:PHE:CD1	1:B:503:ASN:N	2.89	0.40
1:B:242:HIS:NE2	2:B:1002[B]:IOS:O4	2.44	0.40
1:A:513:ILE:HD13	1:A:555:VAL:CG2	2.52	0.40
1:A:464:HIS:CE1	1:A:469:VAL:H	2.40	0.40
1:B:46:VAL:HG22	1:B:73:LYS:HG3	2.04	0.40
1:A:517:SER:O	1:A:520:GLU:N	2.54	0.40
1:B:560:LYS:O	1:B:560:LYS:HG3	2.21	0.40
1:B:359:LYS:CG	1:B:360:CYS:N	2.84	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:A:495:GLU:OE1	1:B:472:ARG:NH2[1_454]	2.13	0.07

5.3 Torsion angles

5.3.1 Protein backbone

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	555/585 (95%)	499 (90%)	44 (8%)	12 (2%)	8	4
1	B	555/585 (95%)	481 (87%)	57 (10%)	17 (3%)	5	2
All	All	1110/1170 (95%)	980 (88%)	101 (9%)	29 (3%)	7	3

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	59	ALA
1	B	312	SER
1	A	57	GLU
1	A	85	GLY
1	A	118	PRO
1	A	511	ALA
1	B	57	GLU
1	B	59	ALA
1	B	85	GLY
1	B	172	ALA
1	B	363	ALA
1	B	442	GLU
1	B	539	ALA
1	A	80	LEU
1	A	172	ALA
1	A	512	ASP
1	B	300	ALA
1	B	512	ASP
1	A	539	ALA
1	B	129	ASP
1	B	163	ALA
1	B	267	ASN
1	B	558	CYS
1	A	150	TYR
1	A	366	PRO

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Mol	Chain	Res	Type
1	B	118	PRO
1	A	77	VAL
1	B	271	ILE
1	B	77	VAL

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	437/511 (86%)	421 (96%)	16 (4%)	41	50
1	B	431/511 (84%)	415 (96%)	16 (4%)	41	50
All	All	868/1022 (85%)	836 (96%)	32 (4%)	41	50

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

Mol	Chain	Res	Type
1	A	37	GLU
1	A	38	ASP
1	A	66	LEU
1	A	84	TYR
1	A	114	ARG
1	A	170	GLN
1	A	208	GLU
1	A	245	CYS
1	A	252	GLU
1	A	334	TYR
1	A	425	GLU
1	A	480	SER
1	A	493	VAL
1	A	503	ASN
1	A	532	LEU
1	A	548	MET
1	B	40	VAL
1	B	84	TYR
1	B	111	ASN

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Mol	Chain	Res	Type
1	B	245	CYS
1	B	252	GLU
1	B	308	ASP
1	B	410	ARG
1	B	419	SER
1	B	425	GLU
1	B	471	ASP
1	B	493	VAL
1	B	503	ASN
1	B	513	ILE
1	B	514	CYS
1	B	532	LEU
1	B	548	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	ASN
1	A	109	ASN
1	A	130	ASN
1	A	196	GLN
1	A	268	GLN
1	A	367	HIS
1	A	385	GLN
1	A	429	ASN
1	A	459	GLN
1	A	464	HIS
1	A	483	ASN
1	A	503	ASN
1	B	9	HIS
1	B	109	ASN
1	B	196	GLN
1	B	247	HIS
1	B	268	GLN
1	B	338	HIS
1	B	367	HIS
1	B	385	GLN
1	B	464	HIS
1	B	483	ASN
1	B	503	ASN
1	B	543	GLN

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 ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	IOS	A	1001	-	15,15,15	2.89	3 (20%)	9,22,22	1.34	1 (11%)
2	IOS	A	1002[A]	-	15,15,15	2.49	5 (33%)	9,22,22	1.26	0
2	IOS	A	1002[B]	-	15,15,15	2.58	5 (33%)	9,22,22	1.20	1 (11%)
2	IOS	B	1001	-	15,15,15	3.03	4 (26%)	9,22,22	1.38	2 (22%)
2	IOS	B	1002[A]	-	15,15,15	2.59	4 (26%)	9,22,22	1.32	2 (22%)
2	IOS	B	1002[B]	-	15,15,15	2.56	5 (33%)	9,22,22	1.29	1 (11%)

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	IOS	A	1001	-	-	0/3/5/5	0/2/2/2
2	IOS	A	1002[A]	-	-	0/3/5/5	0/2/2/2
2	IOS	A	1002[B]	-	-	0/3/5/5	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IOS	B	1001	-	-	0/3/5/5	0/2/2/2
2	IOS	B	1002[A]	-	-	0/3/5/5	0/2/2/2
2	IOS	B	1002[B]	-	-	0/3/5/5	0/2/2/2

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	IOS	O1-S	-10.29	1.44	1.63
2	A	1001	IOS	O1-S	-10.13	1.44	1.63
2	B	1002[A]	IOS	O1-S	-8.24	1.48	1.63
2	A	1002[A]	IOS	O1-S	-7.93	1.49	1.63
2	B	1002[B]	IOS	O1-S	-7.85	1.49	1.63
2	A	1002[B]	IOS	O1-S	-7.78	1.49	1.63
2	A	1001	IOS	C6-C5	2.12	1.43	1.38
2	B	1002[B]	IOS	C5-C4	2.16	1.41	1.36
2	A	1002[A]	IOS	C2-C3	2.23	1.43	1.40
2	A	1002[B]	IOS	C5-C4	2.27	1.41	1.36
2	B	1002[A]	IOS	C5-C4	2.28	1.41	1.36
2	A	1002[A]	IOS	C5-C4	2.34	1.42	1.36
2	A	1002[A]	IOS	C6-C5	2.37	1.44	1.38
2	B	1002[B]	IOS	C2-C3	2.40	1.43	1.40
2	A	1002[B]	IOS	C6-C5	2.48	1.44	1.38
2	B	1001	IOS	C5-C4	2.49	1.42	1.36
2	B	1002[A]	IOS	C6-C5	2.67	1.44	1.38
2	B	1002[B]	IOS	C6-C5	2.71	1.45	1.38
2	B	1001	IOS	C1-N1	2.80	1.42	1.36
2	B	1001	IOS	C6-C5	2.93	1.45	1.38
2	A	1001	IOS	C1-N1	3.02	1.42	1.36
2	A	1002[A]	IOS	C1-N1	3.02	1.42	1.36
2	A	1002[B]	IOS	C2-C3	3.10	1.44	1.40
2	A	1002[B]	IOS	C1-N1	3.35	1.43	1.36
2	B	1002[A]	IOS	C1-N1	3.36	1.43	1.36
2	B	1002[B]	IOS	C1-N1	3.51	1.43	1.36

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	IOS	O4-S-O2	-2.43	99.44	108.56
2	A	1001	IOS	C6-C7-C8	-2.24	116.47	120.06
2	B	1002[A]	IOS	O4-S-O2	-2.18	100.37	108.56
2	A	1002[B]	IOS	O4-S-O2	-2.17	100.38	108.56
2	B	1002[B]	IOS	O4-S-O2	-2.17	100.41	108.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	IOS	C6-C7-C8	-2.15	116.62	120.06
2	B	1002[A]	IOS	C6-C7-C8	-2.02	116.82	120.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1002[A]	IOS	4	0
2	A	1002[B]	IOS	8	0
2	B	1002[A]	IOS	2	0
2	B	1002[B]	IOS	7	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	557/585 (95%)	0.20	32 (5%)	27 30	33, 64, 116, 138	0
1	B	557/585 (95%)	0.18	30 (5%)	29 33	33, 68, 114, 130	0
All	All	1114/1170 (95%)	0.19	62 (5%)	28 31	33, 66, 115, 138	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	508	THR	11.1
1	A	508	THR	6.8
1	A	92	ALA	6.3
1	B	506	THR	5.8
1	B	509	PHE	5.6
1	A	547	VAL	5.5
1	A	83	THR	5.3
1	A	510	HIS	5.1
1	A	561	ALA	4.8
1	B	83	THR	4.7
1	A	513	ILE	4.3
1	B	547	VAL	4.3
1	B	505	GLU	4.3
1	B	171	ALA	4.2
1	A	506	THR	4.1
1	A	511	ALA	4.1
1	A	87	MET	4.1
1	A	551	PHE	4.1
1	A	509	PHE	4.0
1	B	511	ALA	4.0
1	B	551	PHE	3.9
1	A	85	GLY	3.8
1	B	507	PHE	3.7
1	B	555	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	546	ALA	3.6
1	B	554	PHE	3.6
1	B	78	ALA	3.5
1	B	84	TYR	3.4
1	A	505	GLU	3.4
1	A	555	VAL	3.3
1	B	88	ALA	3.3
1	A	502	PHE	3.3
1	B	79	THR	3.2
1	A	91	CYS	3.1
1	B	546	ALA	3.0
1	A	96	PRO	3.0
1	A	516	LEU	3.0
1	A	507	PHE	2.9
1	B	176	ALA	2.9
1	B	558	CYS	2.9
1	B	80	LEU	2.7
1	B	102	PHE	2.7
1	B	168	CYS	2.7
1	A	84	TYR	2.7
1	A	80	LEU	2.6
1	B	515	THR	2.6
1	A	79	THR	2.6
1	B	561	ALA	2.6
1	A	559	CYS	2.6
1	A	81	ARG	2.4
1	A	273	SER	2.4
1	B	510	HIS	2.4
1	A	558	CYS	2.3
1	B	81	ARG	2.3
1	B	82	GLU	2.3
1	A	169	CYS	2.2
1	A	178	LEU	2.2
1	A	512	ASP	2.1
1	B	74	LEU	2.1
1	B	548	MET	2.1
1	B	460	LEU	2.1
1	A	554	PHE	2.0

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	IOS	A	1002[B]	14/14	0.70	0.46	8.38	79,80,81,81	14
2	IOS	A	1002[A]	14/14	0.70	0.46	5.25	108,109,111,111	14
2	IOS	B	1002[B]	14/14	0.82	0.37	4.03	74,74,76,77	14
2	IOS	B	1002[A]	14/14	0.82	0.37	3.27	108,109,110,110	14
2	IOS	A	1001	14/14	0.99	0.18	1.13	40,43,50,51	0
2	IOS	B	1001	14/14	0.98	0.15	0.35	49,51,55,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.