



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:52 AM GMT

PDB ID : 2J4C
Title : STRUCTURE OF HUMAN BUTYRYLCHOLINESTERASE IN COMPLEX WITH 10MM HGCL2
Authors : Colletier, J.P.; Frasco, M.F.; Carvalho, F.; Guilhermino, L.; Stojan, J.; Fournier, D.; Weik, M.
Deposited on : 2006-08-28
Resolution : 2.75 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

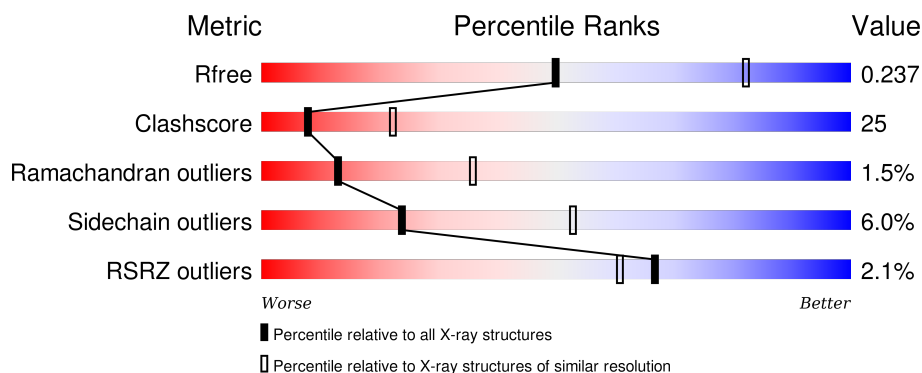
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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	529	

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
3	NAG	A	1538	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	CL	A	1539	-	-	X	-
6	SO4	A	1549	-	-	X	-
7	GOL	A	1544	-	-	-	X
7	GOL	A	1545	-	-	-	X
7	GOL	A	1546	-	-	-	X
8	BUA	A	1547	-	-	X	X
9	MES	A	1548	-	-	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 4712 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 CHOLINESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	524	Total	C	N	O	S	0	0	0
			4176	2698	702	760	16			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	GLN	ASN	ENGINEERED MUTATION	UNP P06276
A	455	GLN	ASN	ENGINEERED MUTATION	UNP P06276
A	481	GLN	ASN	ENGINEERED MUTATION	UNP P06276
A	486	ALA	ASN	ENGINEERED MUTATION	UNP P06276

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	3	Total	C	N	O	0	0
			38	22	2	14		
2	A	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

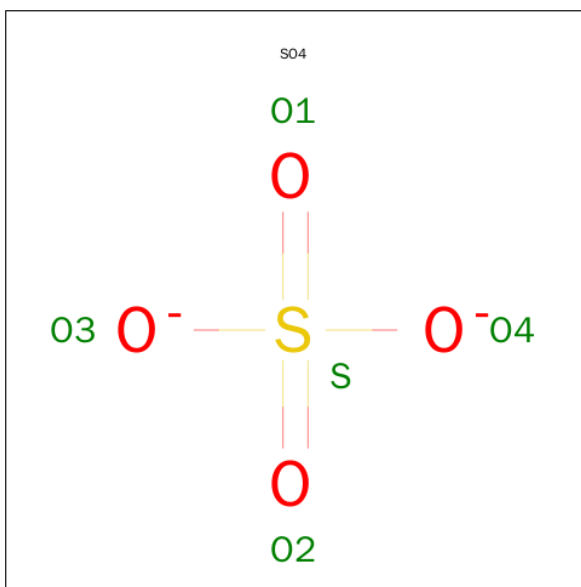
- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Cl	0	0
			3	3		

- Molecule 5 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

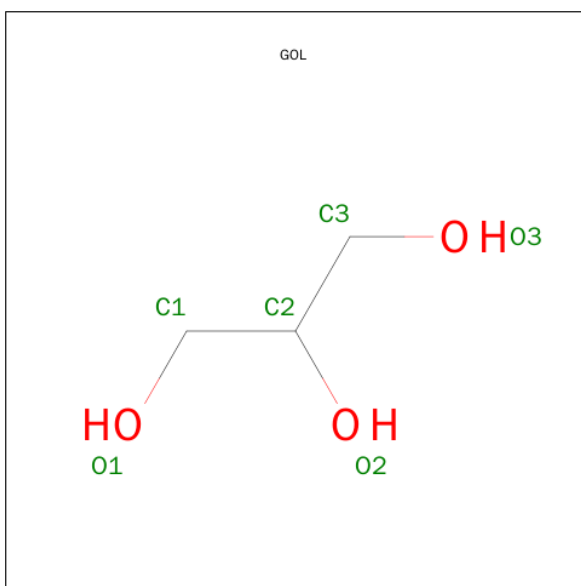
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	Hg	0	0
			4	4		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



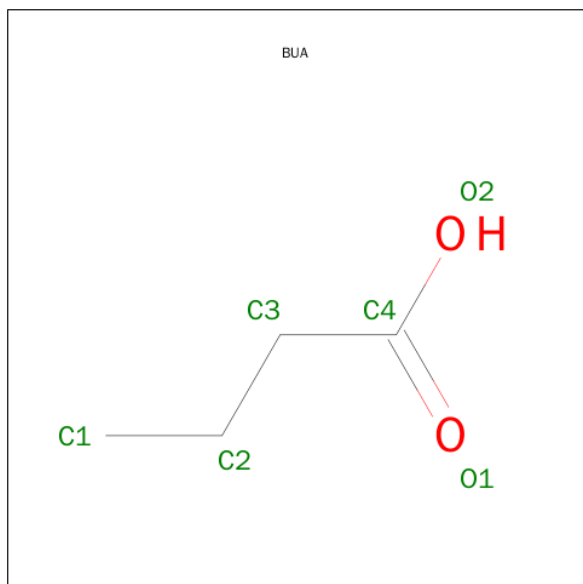
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

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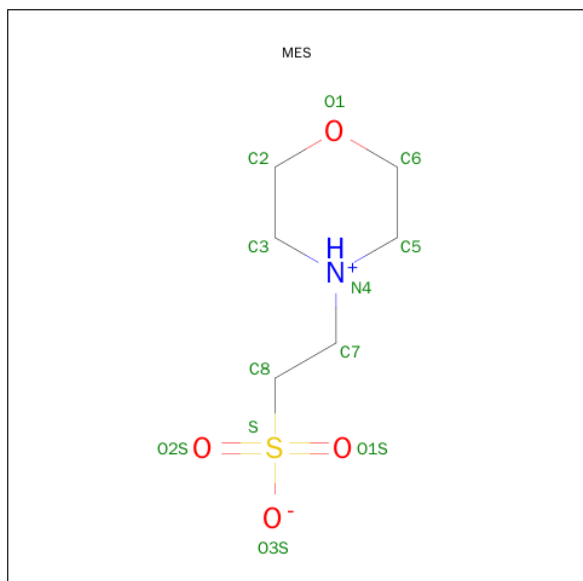
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is BUTANOIC ACID (three-letter code: BUA) (formula: $C_4H_8O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			6	4	2		

- Molecule 9 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

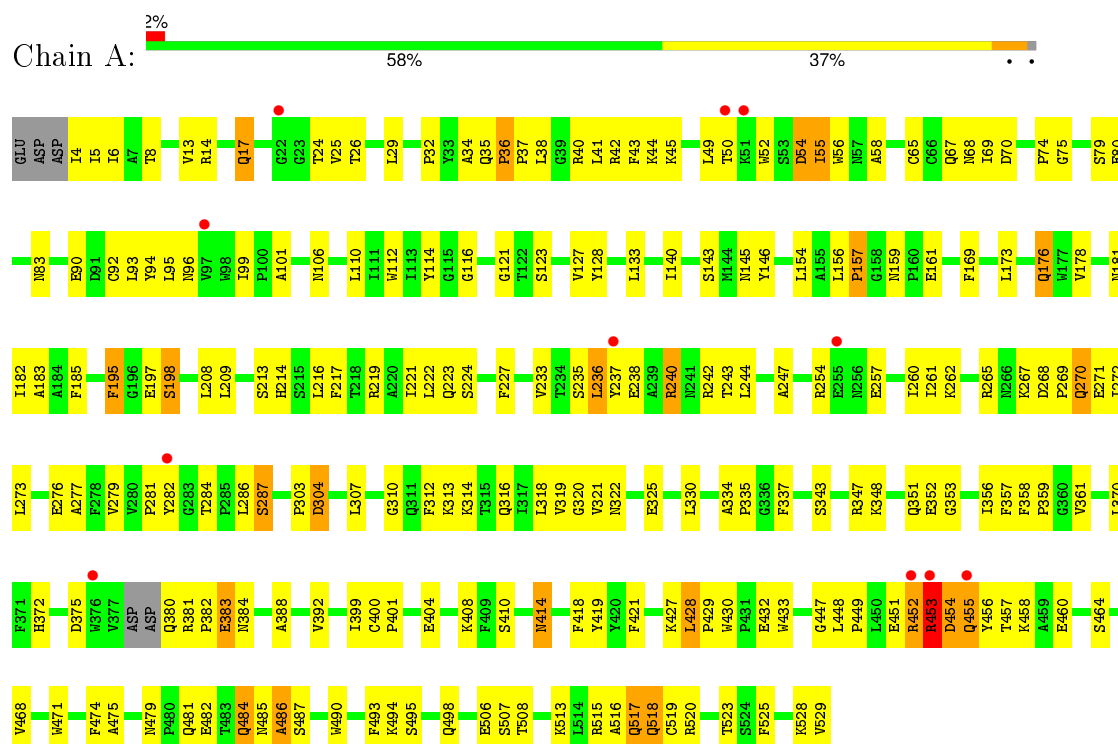
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	365	Total	O	0	0
			365	365		

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: CHOLINESTERASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	153.76 Å 153.76 Å 128.58 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.39 – 2.75 19.39 – 2.75	Depositor EDS
% Data completeness (in resolution range)	93.9 (19.39-2.75) 97.7 (19.39-2.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 2.75 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.184 , 0.231 0.195 , 0.237	Depositor DCC
R_{free} test set	989 reflections (5.44%)	DCC
Wilson B-factor (Å ²)	51.7	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 57.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 37308 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4712	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% 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: GOL, NAG, CL, FUC, MES, BUA, HG, CSS, 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	0.69	0/4286	0.74	1/5815 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	453	ARG	N-CA-C	-5.45	96.29	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4176	0	4086	207	0
2	A	76	0	68	2	0
3	A	42	0	39	2	0
4	A	3	0	0	4	0
5	A	4	0	0	0	0
6	A	10	0	0	2	0
7	A	18	0	24	5	0
8	A	6	0	8	7	0
9	A	12	0	13	2	0
10	A	365	0	0	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4712	0	4238	214	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 25.

All (214) 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:451:GLU:HG2	1:A:454:ASP:OD1	1.55	1.04
1:A:198:SER:OG	8:A:1547:BUA:C4	2.16	0.92
1:A:304:ASP:HB3	10:A:2077:HOH:O	1.69	0.90
1:A:518:GLN:H	1:A:518:GLN:NE2	1.71	0.88
1:A:286:LEU:HD23	8:A:1547:BUA:H22	1.56	0.87
1:A:242:ARG:HH22	7:A:1545:GOL:H2	1.42	0.83
1:A:518:GLN:H	1:A:518:GLN:HE21	1.25	0.83
1:A:372:HIS:HB2	6:A:1549:SO4:O4	1.79	0.82
7:A:1546:GOL:H31	10:A:2358:HOH:O	1.79	0.80
1:A:154:LEU:HD11	1:A:243:THR:HG23	1.64	0.79
1:A:96:ASN:HD22	1:A:143:SER:HB2	1.48	0.78
1:A:372:HIS:ND1	6:A:1549:SO4:O1	2.17	0.78
1:A:209:LEU:HD23	1:A:312:PHE:HB3	1.70	0.74
1:A:375:ASP:HB3	10:A:2245:HOH:O	1.88	0.74
1:A:348:LYS:HE3	1:A:352:GLU:OE2	1.87	0.74
1:A:381:ARG:HB3	1:A:384:ASN:ND2	2.03	0.73
4:A:3001:CL:CL	10:A:2237:HOH:O	2.43	0.72
1:A:54:ASP:HB3	10:A:2045:HOH:O	1.90	0.71
1:A:286:LEU:CD2	8:A:1547:BUA:H22	2.20	0.71
1:A:453:ARG:N	1:A:453:ARG:CD	2.53	0.71
1:A:330:LEU:O	1:A:334:ALA:HB3	1.92	0.69
1:A:453:ARG:O	1:A:455:GLN:N	2.25	0.69
1:A:198:SER:HG	8:A:1547:BUA:C4	2.05	0.69
8:A:1547:BUA:H11	10:A:2198:HOH:O	1.92	0.68
1:A:106:ASN:HB2	3:A:1536:NAG:H83	1.75	0.67
1:A:451:GLU:CG	1:A:454:ASP:OD1	2.40	0.66
1:A:358:PHE:HB3	1:A:361:VAL:HG21	1.78	0.66
1:A:515:ARG:NH2	9:A:1548:MES:H72	2.11	0.66
1:A:26:THR:HB	1:A:99:ILE:HG12	1.78	0.66
1:A:69:ILE:HG23	1:A:83:ASN:OD1	1.95	0.66
1:A:37:PRO:HG3	1:A:93:LEU:HD11	1.78	0.65
1:A:156:LEU:HD13	1:A:243:THR:HG21	1.79	0.65
1:A:262:LYS:HG2	10:A:2186:HOH:O	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:LEU:HD13	1:A:430:TRP:H	1.63	0.63
1:A:357:PHE:O	1:A:359:PRO:HD2	1.98	0.63
1:A:427:LYS:HE3	1:A:455:GLN:HB2	1.79	0.63
1:A:257:GLU:HG2	10:A:2181:HOH:O	1.99	0.62
1:A:516:ALA:O	1:A:520:ARG:HG3	1.98	0.62
1:A:123:SER:HB2	1:A:143:SER:HB3	1.79	0.62
1:A:388:ALA:O	1:A:392:VAL:HG23	2.01	0.61
1:A:34:ALA:HB2	1:A:173:LEU:HD23	1.81	0.61
1:A:404:GLU:HG2	10:A:2166:HOH:O	2.01	0.61
1:A:455:GLN:N	1:A:455:GLN:OE1	2.34	0.60
1:A:240:ARG:HG2	1:A:240:ARG:HH11	1.65	0.60
1:A:383:GLU:CD	1:A:383:GLU:H	2.04	0.60
1:A:284:THR:H	1:A:287:SER:HB2	1.64	0.60
1:A:5:ILE:HD13	1:A:55:ILE:HD13	1.84	0.60
7:A:1546:GOL:C3	8:A:1547:BUA:O1	2.50	0.59
1:A:380:GLN:HB2	10:A:2252:HOH:O	2.02	0.59
1:A:518:GLN:HE21	1:A:518:GLN:N	1.97	0.59
1:A:493:PHE:CZ	1:A:498:GLN:HB3	2.37	0.59
1:A:227:PHE:HZ	1:A:307:LEU:HB2	1.68	0.59
1:A:270:GLN:HB2	10:A:2187:HOH:O	2.02	0.59
1:A:381:ARG:HB3	1:A:384:ASN:HD21	1.68	0.58
1:A:381:ARG:HG3	1:A:382:PRO:HD2	1.85	0.58
1:A:508:THR:HG22	4:A:1539:CL:CL	2.40	0.58
1:A:213:SER:HA	1:A:216:LEU:HD12	1.85	0.58
1:A:58:ALA:HA	10:A:2020:HOH:O	2.04	0.58
1:A:262:LYS:HG3	1:A:265:ARG:NH1	2.19	0.57
1:A:319:VAL:O	1:A:418:PHE:HA	2.04	0.57
1:A:318:LEU:HD21	1:A:490:TRP:CH2	2.39	0.57
1:A:453:ARG:N	1:A:453:ARG:HD2	2.19	0.57
1:A:238:GLU:O	1:A:242:ARG:HG3	2.04	0.57
1:A:452:ARG:HH21	1:A:458:LYS:HD3	1.69	0.57
1:A:404:GLU:OE1	1:A:523:THR:HG21	2.04	0.57
1:A:37:PRO:HG3	1:A:93:LEU:CD1	2.34	0.57
1:A:318:LEU:HD21	1:A:490:TRP:HH2	1.69	0.57
1:A:370:LEU:HD23	1:A:370:LEU:C	2.25	0.57
1:A:29:LEU:O	1:A:58:ALA:HB1	2.04	0.56
1:A:244:LEU:O	1:A:247:ALA:HB3	2.05	0.56
1:A:357:PHE:C	1:A:359:PRO:CD	2.74	0.56
1:A:198:SER:HA	1:A:224:SER:O	2.04	0.56
1:A:17:GLN:HE21	1:A:17:GLN:HA	1.70	0.56
1:A:237:TYR:HB2	10:A:2169:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:LYS:CE	1:A:454:ASP:O	2.53	0.56
1:A:427:LYS:HE3	1:A:454:ASP:O	2.06	0.56
1:A:240:ARG:HD2	10:A:2171:HOH:O	2.06	0.55
8:A:1547:BUA:C1	10:A:2360:HOH:O	2.53	0.55
1:A:183:ALA:HB2	10:A:2145:HOH:O	2.05	0.55
1:A:310:GLY:HA2	1:A:312:PHE:CE2	2.42	0.55
1:A:75:GLY:H	1:A:80:GLU:CD	2.11	0.55
1:A:508:THR:CG2	4:A:1539:CL:CL	2.92	0.55
1:A:382:PRO:HG2	1:A:383:GLU:OE1	2.07	0.54
1:A:453:ARG:N	1:A:453:ARG:HD3	2.24	0.53
1:A:358:PHE:HB3	1:A:361:VAL:CG2	2.38	0.53
1:A:65:CYS:O	1:A:67:GLN:HG2	2.07	0.53
1:A:214:HIS:O	1:A:313:LYS:HE2	2.09	0.53
1:A:314:LYS:HB3	1:A:414:ASN:HD21	1.74	0.53
1:A:56:TRP:CD1	1:A:56:TRP:C	2.81	0.53
1:A:41:LEU:HB3	10:A:2032:HOH:O	2.08	0.53
1:A:498:GLN:OE1	1:A:513:LYS:HG2	2.08	0.53
1:A:320:GLY:HA3	1:A:419:TYR:CE1	2.44	0.52
1:A:432:GLU:HG2	10:A:2279:HOH:O	2.10	0.52
1:A:494:LYS:HD2	10:A:2310:HOH:O	2.07	0.52
1:A:235:SER:OG	1:A:238:GLU:HG3	2.10	0.52
1:A:262:LYS:HG3	1:A:265:ARG:HH12	1.74	0.51
1:A:24:THR:O	1:A:101:ALA:HB3	2.11	0.51
1:A:353:GLY:O	1:A:356:ILE:HB	2.11	0.51
1:A:281:PRO:HG2	1:A:282:TYR:H	1.75	0.51
1:A:112:TRP:HA	1:A:195:PHE:O	2.11	0.51
1:A:68:ASN:OD1	1:A:273:LEU:HB3	2.10	0.51
1:A:242:ARG:HH22	7:A:1545:GOL:C2	2.19	0.50
1:A:42:ARG:O	1:A:43:PHE:HB2	2.12	0.50
1:A:4:ILE:HD13	1:A:17:GLN:HG2	1.91	0.50
1:A:197:GLU:HA	1:A:223:GLN:O	2.11	0.50
1:A:525:PHE:HA	1:A:528:LYS:HE2	1.94	0.50
1:A:159:ASN:HB2	10:A:2182:HOH:O	2.11	0.50
2:A:1533:NAG:H82	10:A:2345:HOH:O	2.11	0.50
1:A:457:THR:OG1	1:A:460:GLU:HG3	2.12	0.50
1:A:357:PHE:C	1:A:359:PRO:HD2	2.33	0.49
1:A:32:PRO:HB3	1:A:94:TYR:CE2	2.47	0.49
1:A:337:PHE:HA	1:A:343:SER:OG	2.12	0.49
1:A:50:THR:HB	10:A:2040:HOH:O	2.11	0.49
1:A:156:LEU:HD22	1:A:261:ILE:HD11	1.95	0.49
1:A:358:PHE:N	1:A:359:PRO:CD	2.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:TRP:HZ2	10:A:2018:HOH:O	1.95	0.49
1:A:176:GLN:HG3	10:A:2140:HOH:O	2.12	0.49
1:A:428:LEU:CD1	1:A:430:TRP:HB2	2.43	0.48
1:A:276:GLU:O	1:A:279:VAL:HG22	2.13	0.48
1:A:42:ARG:HG2	1:A:43:PHE:CD2	2.49	0.48
1:A:112:TRP:HB2	1:A:195:PHE:CE1	2.48	0.48
1:A:13:VAL:HG12	1:A:52:TRP:HZ2	1.79	0.48
1:A:348:LYS:O	1:A:352:GLU:HG2	2.14	0.48
1:A:6:ILE:HD12	1:A:185:PHE:CD2	2.49	0.48
1:A:114:TYR:CE1	1:A:146:TYR:CE2	3.01	0.48
1:A:195:PHE:CB	1:A:221:ILE:HB	2.44	0.47
1:A:209:LEU:CD2	1:A:312:PHE:HB3	2.42	0.47
1:A:219:ARG:HD3	1:A:316:GLN:OE1	2.14	0.47
1:A:227:PHE:CD2	1:A:303:PRO:HB2	2.50	0.47
1:A:453:ARG:H	1:A:453:ARG:HD3	1.80	0.47
1:A:209:LEU:HD11	1:A:227:PHE:CE1	2.50	0.47
1:A:95:LEU:HD12	1:A:95:LEU:C	2.35	0.47
1:A:242:ARG:NH2	7:A:1545:GOL:H2	2.20	0.47
1:A:321:VAL:HG11	1:A:399:ILE:HA	1.96	0.47
1:A:508:THR:N	10:A:2320:HOH:O	2.48	0.46
1:A:449:PRO:HA	1:A:456:TYR:CD1	2.50	0.46
1:A:400:CYS:HB2	1:A:401:PRO:HD3	1.97	0.46
1:A:70:ASP:OD1	1:A:79:SER:HB2	2.16	0.46
1:A:8:THR:HB	1:A:181:ASN:OD1	2.16	0.46
1:A:99:ILE:HG21	1:A:185:PHE:HB3	1.98	0.46
1:A:36:PRO:HA	1:A:37:PRO:HD3	1.84	0.46
1:A:508:THR:HB	10:A:2321:HOH:O	2.15	0.45
1:A:157:PRO:HG2	1:A:236:LEU:CD2	2.46	0.45
1:A:479:ASN:OD1	1:A:481:GLN:HG2	2.15	0.45
1:A:428:LEU:HD11	1:A:430:TRP:HB2	1.97	0.45
1:A:519:CYS:HB3	10:A:2336:HOH:O	2.16	0.45
1:A:433:TRP:HE3	1:A:433:TRP:O	1.99	0.45
1:A:195:PHE:HB2	1:A:221:ILE:HB	1.99	0.45
1:A:357:PHE:C	1:A:359:PRO:HD3	2.37	0.45
1:A:372:HIS:CE1	1:A:517:GLN:OE1	2.69	0.45
1:A:217:PHE:O	1:A:313:LYS:HE3	2.17	0.44
1:A:240:ARG:HG2	1:A:240:ARG:NH1	2.32	0.44
1:A:45:LYS:HD2	1:A:169:PHE:CD1	2.52	0.44
1:A:106:ASN:CB	3:A:1536:NAG:H83	2.45	0.44
1:A:222:LEU:N	1:A:222:LEU:HD12	2.32	0.44
1:A:322:ASN:ND2	1:A:421:PHE:HB3	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ARG:HG2	10:A:2031:HOH:O	2.17	0.44
1:A:123:SER:HB2	1:A:143:SER:CB	2.47	0.44
1:A:347:ARG:O	1:A:351:GLN:HG3	2.17	0.44
1:A:209:LEU:HD11	1:A:227:PHE:CZ	2.53	0.44
1:A:284:THR:HG22	1:A:356:ILE:O	2.17	0.44
1:A:272:ILE:O	1:A:276:GLU:HG3	2.18	0.43
1:A:277:ALA:HB2	10:A:2058:HOH:O	2.18	0.43
1:A:221:ILE:C	1:A:222:LEU:HD12	2.38	0.43
1:A:518:GLN:NE2	1:A:518:GLN:N	2.53	0.43
1:A:37:PRO:HD2	1:A:90:GLU:O	2.18	0.43
1:A:49:LEU:HD12	1:A:50:THR:H	1.84	0.43
1:A:529:VAL:HB	10:A:2337:HOH:O	2.18	0.43
1:A:178:VAL:HG13	1:A:182:ILE:HB	2.01	0.43
1:A:17:GLN:NE2	1:A:26:THR:HG23	2.33	0.43
1:A:74:PRO:HA	1:A:80:GLU:OE1	2.19	0.43
1:A:233:VAL:HG21	1:A:303:PRO:HG3	2.00	0.43
1:A:92:CYS:O	1:A:145:ASN:HB2	2.19	0.43
1:A:127:VAL:HG23	1:A:128:TYR:CD1	2.54	0.43
1:A:484:GLN:O	1:A:486:ALA:N	2.51	0.43
1:A:5:ILE:HD11	1:A:14:ARG:NH1	2.34	0.43
1:A:318:LEU:HD12	1:A:474:PHE:CG	2.55	0.42
1:A:254:ARG:H	1:A:260:ILE:HG12	1.83	0.42
1:A:334:ALA:HA	1:A:335:PRO:HD2	1.87	0.42
1:A:448:LEU:N	1:A:449:PRO:CD	2.82	0.42
1:A:133:LEU:HD23	1:A:468:VAL:HG13	2.01	0.42
1:A:43:PHE:CE1	1:A:154:LEU:HD22	2.54	0.42
1:A:17:GLN:NE2	1:A:17:GLN:HA	2.33	0.42
1:A:35:GLN:O	1:A:36:PRO:C	2.54	0.42
1:A:322:ASN:HD22	1:A:421:PHE:HB3	1.84	0.42
1:A:140:ILE:HG21	1:A:182:ILE:HD11	2.00	0.42
1:A:183:ALA:HB3	10:A:2144:HOH:O	2.19	0.42
1:A:24:THR:CG2	1:A:25:VAL:N	2.82	0.42
2:A:1533:NAG:O6	2:A:1535:FUC:H4	2.20	0.42
1:A:44:LYS:HG3	10:A:2035:HOH:O	2.19	0.42
1:A:80:GLU:HA	1:A:83:ASN:ND2	2.35	0.41
1:A:208:LEU:HD23	1:A:217:PHE:CZ	2.54	0.41
1:A:447:GLY:HA2	1:A:464:SER:OG	2.20	0.41
1:A:110:LEU:HD11	1:A:475:ALA:CB	2.50	0.41
1:A:348:LYS:HA	1:A:348:LYS:HD2	1.79	0.41
1:A:321:VAL:HG21	1:A:399:ILE:HG12	2.01	0.41
1:A:123:SER:CB	1:A:143:SER:HB3	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:507:SER:HB2	10:A:2318:HOH:O	2.21	0.41
1:A:453:ARG:NH1	1:A:453:ARG:HG3	2.36	0.41
1:A:427:LYS:NZ	1:A:454:ASP:O	2.54	0.41
1:A:508:THR:HG21	4:A:1539:CL:CL	2.58	0.41
1:A:485:ASN:N	1:A:485:ASN:OD1	2.53	0.41
1:A:428:LEU:HA	1:A:429:PRO:HD2	1.83	0.41
1:A:404:GLU:OE2	1:A:408:LYS:HE2	2.20	0.41
1:A:67:GLN:N	1:A:273:LEU:HD11	2.35	0.41
1:A:114:TYR:HE2	1:A:145:ASN:HA	1.86	0.41
1:A:515:ARG:HH22	9:A:1548:MES:H72	1.83	0.41
1:A:322:ASN:O	1:A:325:GLU:HG2	2.21	0.41
1:A:178:VAL:HB	10:A:2138:HOH:O	2.21	0.40
1:A:55:ILE:HG13	1:A:55:ILE:H	1.71	0.40
1:A:183:ALA:CB	10:A:2145:HOH:O	2.67	0.40
1:A:267:LYS:HD2	1:A:271:GLU:OE2	2.21	0.40
1:A:121:GLY:HA2	10:A:2072:HOH:O	2.21	0.40
1:A:159:ASN:OD1	1:A:161:GLU:HG2	2.21	0.40

There are no symmetry-related clashes.

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	519/529 (98%)	458 (88%)	53 (10%)	8 (2%)	13	36

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	454	ASP
1	A	486	ALA
1	A	506	GLU
1	A	55	ILE

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Mol	Chain	Res	Type
1	A	453	ARG
1	A	54	ASP
1	A	116	GLY
1	A	157	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	447/452 (99%)	420 (94%)	27 (6%)	24 53

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	36	PRO
1	A	38	LEU
1	A	176	GLN
1	A	195	PHE
1	A	198	SER
1	A	236	LEU
1	A	240	ARG
1	A	268	ASP
1	A	269	PRO
1	A	270	GLN
1	A	287	SER
1	A	304	ASP
1	A	383	GLU
1	A	410	SER
1	A	414	ASN
1	A	428	LEU
1	A	452	ARG
1	A	453	ARG
1	A	455	GLN
1	A	471	TRP
1	A	482	GLU

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Mol	Chain	Res	Type
1	A	484	GLN
1	A	487	SER
1	A	495	SER
1	A	517	GLN
1	A	518	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	47	GLN
1	A	172	GLN
1	A	176	GLN
1	A	214	HIS
1	A	322	ASN
1	A	384	ASN
1	A	518	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	$\# Z > 2$
1	CSS	A	66	1	4,6,7	0.46	0	3,6,8	0.84	0

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	CSS	A	66	1	-	0/1/5/7	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

6 carbohydrates 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	NAG	A	1530	1,2	14,14,15	0.49	0	15,19,21	0.89	1 (6%)
2	NAG	A	1531	2	14,14,15	0.54	0	15,19,21	0.88	1 (6%)
2	FUC	A	1532	2	10,10,11	0.48	0	14,14,16	0.50	0
2	NAG	A	1533	1,2	14,14,15	0.67	0	15,19,21	0.62	0
2	NAG	A	1534	2	14,14,15	0.38	0	15,19,21	0.80	1 (6%)
2	FUC	A	1535	2	10,10,11	0.95	1 (10%)	14,14,16	0.97	1 (7%)

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	NAG	A	1530	1,2	-	1/6/23/26	0/1/1/1
2	NAG	A	1531	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FUC	A	1532	2	-	0/0/17/20	0/1/1/1
2	NAG	A	1533	1,2	-	1/6/23/26	0/1/1/1
2	NAG	A	1534	2	-	0/6/23/26	0/1/1/1
2	FUC	A	1535	2	-	0/0/17/20	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1535	FUC	C2-C3	2.18	1.55	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1534	NAG	C2-N2-C7	-2.47	119.86	123.04
2	A	1531	NAG	C2-N2-C7	-2.47	119.87	123.04
2	A	1530	NAG	C2-N2-C7	-2.38	119.99	123.04
2	A	1535	FUC	C3-C4-C5	2.29	113.57	109.72

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1530	NAG	O7-C7-N2-C2
2	A	1533	NAG	O7-C7-N2-C2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1533	NAG	2	0
2	A	1535	FUC	1	0

5.6 Ligand geometry ⓘ

Of 17 ligands modelled in this entry, 7 are monoatomic - leaving 10 for Mogul analysis.

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	$\# Z > 2$
3	NAG	A	1536	1	14,14,15	0.70	0	15,19,21	0.78	1 (6%)
3	NAG	A	1537	1	14,14,15	0.66	0	15,19,21	0.79	1 (6%)
3	NAG	A	1538	1	14,14,15	0.71	0	15,19,21	0.63	0
6	SO4	A	1543	-	4,4,4	0.34	0	6,6,6	0.14	0
7	GOL	A	1544	-	5,5,5	0.26	0	5,5,5	0.30	0
7	GOL	A	1545	-	5,5,5	0.30	0	5,5,5	0.20	0
7	GOL	A	1546	-	5,5,5	0.26	0	5,5,5	0.51	0
8	BUA	A	1547	-	2,5,5	0.58	0	1,5,5	1.08	0
9	MES	A	1548	-	11,12,12	1.13	0	14,16,16	1.42	1 (7%)
6	SO4	A	1549	5	4,4,4	0.76	0	6,6,6	0.15	0

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
3	NAG	A	1536	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1537	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1538	1	-	0/6/23/26	0/1/1/1
6	SO4	A	1543	-	-	0/0/0/0	0/0/0/0
7	GOL	A	1544	-	-	0/4/4/4	0/0/0/0
7	GOL	A	1545	-	-	0/4/4/4	0/0/0/0
7	GOL	A	1546	-	-	0/4/4/4	0/0/0/0
8	BUA	A	1547	-	-	0/1/3/3	0/0/0/0
9	MES	A	1548	-	-	0/6/14/14	0/1/1/1
6	SO4	A	1549	5	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1537	NAG	C2-N2-C7	-2.35	120.02	123.04
3	A	1536	NAG	C2-N2-C7	-2.33	120.04	123.04
9	A	1548	MES	C5-N4-C3	3.93	117.41	108.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1536	NAG	2	0
7	A	1545	GOL	3	0
7	A	1546	GOL	2	0
8	A	1547	BUA	7	0
9	A	1548	MES	2	0
6	A	1549	SO4	2	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 [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	523/529 (98%)	-0.26	11 (2%) 67 61	29, 59, 95, 147	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	282	TYR	4.6
1	A	51	LYS	4.6
1	A	452	ARG	4.0
1	A	255	GLU	3.7
1	A	453	ARG	2.9
1	A	237	TYR	2.9
1	A	455	GLN	2.7
1	A	376	TRP	2.6
1	A	97	VAL	2.5
1	A	22	GLY	2.4
1	A	50	THR	2.3

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. 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(Å ²)	Q<0.9
1	CSS	A	66	7/8	0.95	0.11	-	61,61,61,61	0

6.3 Carbohydrates ⓘ

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(Å ²)	Q<0.9
2	NAG	A	1533	14/15	0.94	0.24	1.97	72,72,72,72	0
2	FUC	A	1532	10/11	0.96	0.21	-0.13	101,101,101,101	0
2	FUC	A	1535	10/11	0.86	0.38	-	105,105,105,105	0
2	NAG	A	1531	14/15	0.85	0.40	-	108,108,108,108	0
2	NAG	A	1530	14/15	0.92	0.19	-	125,125,125,125	0
2	NAG	A	1534	14/15	0.91	0.26	-	101,101,101,101	0

6.4 Ligands ⓘ

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(Å ²)	Q<0.9
7	GOL	A	1546	6/6	0.75	0.31	5.27	61,61,61,61	0
7	GOL	A	1545	6/6	0.82	0.31	4.44	61,61,61,61	0
7	GOL	A	1544	6/6	0.89	0.34	4.00	61,61,61,61	0
3	NAG	A	1538	14/15	0.83	0.29	3.46	101,101,101,101	0
8	BUA	A	1547	6/6	0.87	0.21	2.93	72,72,72,72	0
9	MES	A	1548	12/12	0.86	0.28	2.34	141,141,141,141	0
6	SO4	A	1549	5/5	0.93	0.14	-0.87	103,103,103,103	0
5	HG	A	1541	1/1	0.95	0.05	-1.85	112,112,112,112	1
4	CL	A	3001	1/1	0.89	0.08	-	83,83,83,83	0
5	HG	A	3003	1/1	0.86	0.10	-	129,129,129,129	1
5	HG	A	1542	1/1	0.89	0.20	-	209,209,209,209	1
4	CL	A	3002	1/1	0.95	0.11	-	100,100,100,100	0
4	CL	A	1539	1/1	0.94	0.06	-	78,78,78,78	0
5	HG	A	1540	1/1	0.95	0.20	-	104,104,104,104	1
3	NAG	A	1537	14/15	0.76	0.30	-	111,111,111,111	0
3	NAG	A	1536	14/15	0.91	0.30	-	80,80,80,80	0
6	SO4	A	1543	5/5	0.96	0.27	-	106,106,106,106	0

6.5 Other polymers

There are no such residues in this entry.