



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:52 AM GMT

PDB ID : 2C10
Title : THE STRUCTURE OF A TRUNCATED, SOLUBLE VERSION OF SEMIC
ARBAZIDE-SENSITIVE AMINE OXIDASE
Authors : Jakobsson, E.; Kleywegt, G.J.
Deposited on : 2005-09-09
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
<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

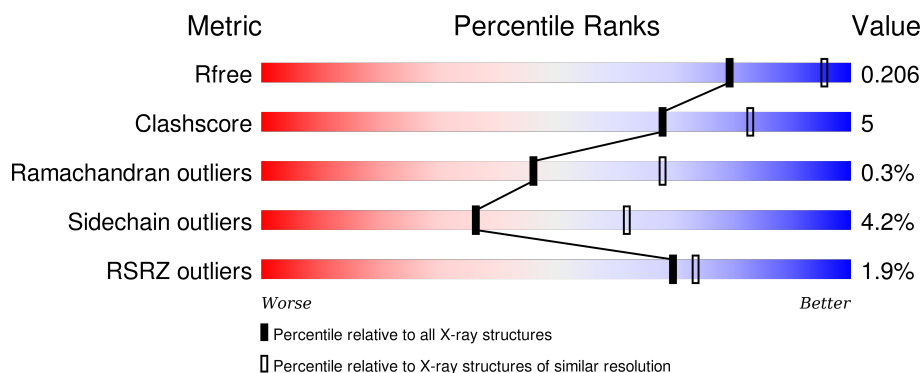
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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 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	735	<div> <div>2%</div> <div>81% 13% . .</div> </div>
1	B	735	<div> <div>%</div> <div>84% 11% . .</div> </div>
1	C	735	<div> <div>2%</div> <div>84% 11% . .</div> </div>
1	D	735	<div> <div>2%</div> <div>85% 10% . .</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
4	NAG	A	1767	-	-	-	X
4	NAG	A	1770	-	-	-	X
4	NAG	C	1770	-	-	-	X

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 23470 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 MEMBRANE COPPER AMINE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	706	Total	C	N	O	S	0	0	1
			5563	3571	960	1011	21			
1	B	709	Total	C	N	O	S	0	0	1
			5587	3587	965	1014	21			
1	C	706	Total	C	N	O	S	0	0	1
			5563	3571	960	1011	21			
1	D	709	Total	C	N	O	S	0	0	1
			5587	3587	965	1014	21			

- 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
			39	22	2	15		
2	B	3	Total	C	N	O	0	0
			39	22	2	15		
2	C	3	Total	C	N	O	0	0
			39	22	2	15		

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

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

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	0
			28	16	2	10		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	2	Total	C	N	O	0	0
			28	16	2	10		
4	C	2	Total	C	N	O	0	0
			28	16	2	10		
4	D	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total 2	Ca 2	0	0
6	A	2	Total 2	Ca 2	0	0
6	D	2	Total 2	Ca 2	0	0
6	C	2	Total 2	Ca 2	0	0

- Molecule 7 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total 1	Cu 1	0	0
7	A	1	Total 1	Cu 1	0	0
7	D	1	Total 1	Cu 1	0	0
7	C	1	Total 1	Cu 1	0	0

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	2	Total 2	Cl 2	0	0
8	A	2	Total 2	Cl 2	0	0
8	D	2	Total 2	Cl 2	0	0
8	C	2	Total 2	Cl 2	0	0

- Molecule 9 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	B	5	Total 60	C 34	N 2	O 24	0	0
9	D	5	Total 60	C 34	N 2	O 24	0	0

- Molecule 10 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	D	4	Total	C	N	O	0	0
			50	28	2	20		

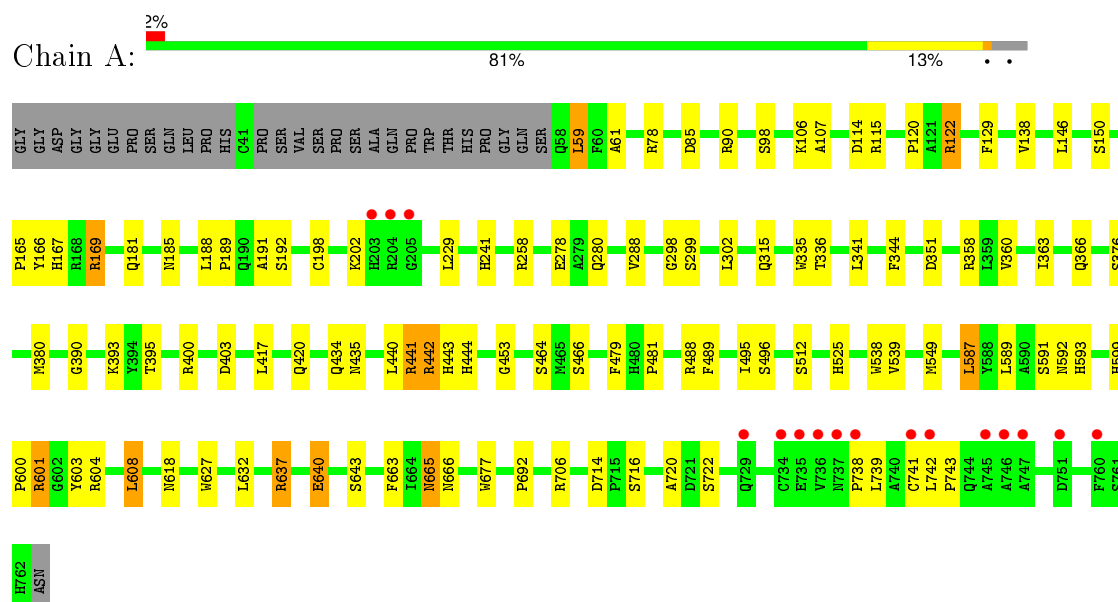
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	141	Total	O	0	0
			141	141		
11	B	139	Total	O	0	0
			139	139		
11	C	141	Total	O	0	0
			141	141		
11	D	142	Total	O	0	0
			142	142		

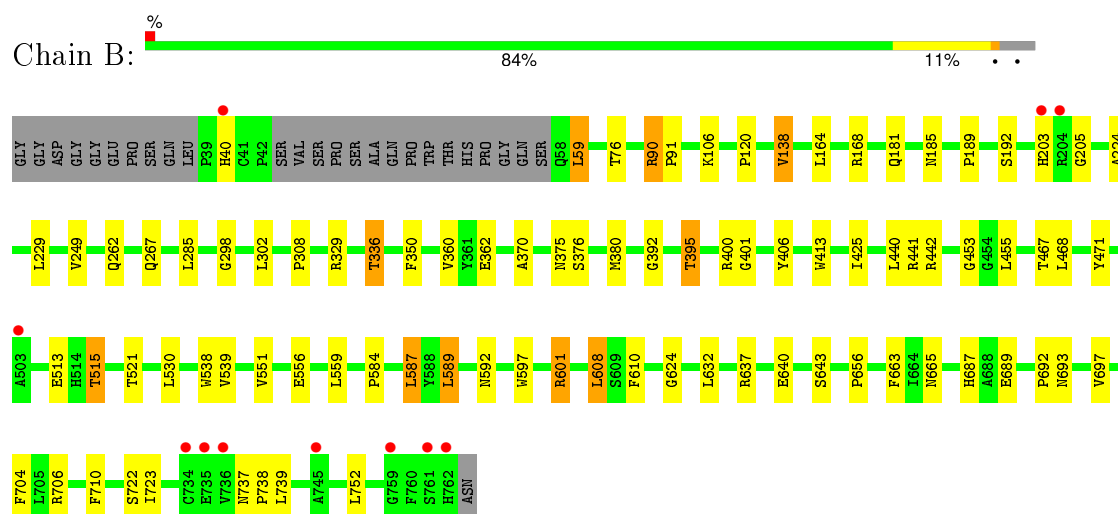
3 Residue-property plots

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.

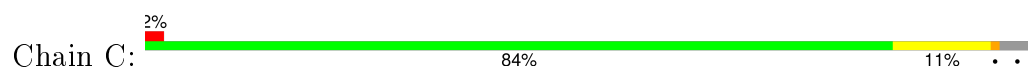
• Molecule 1: MEMBRANE COPPER AMINE OXIDASE

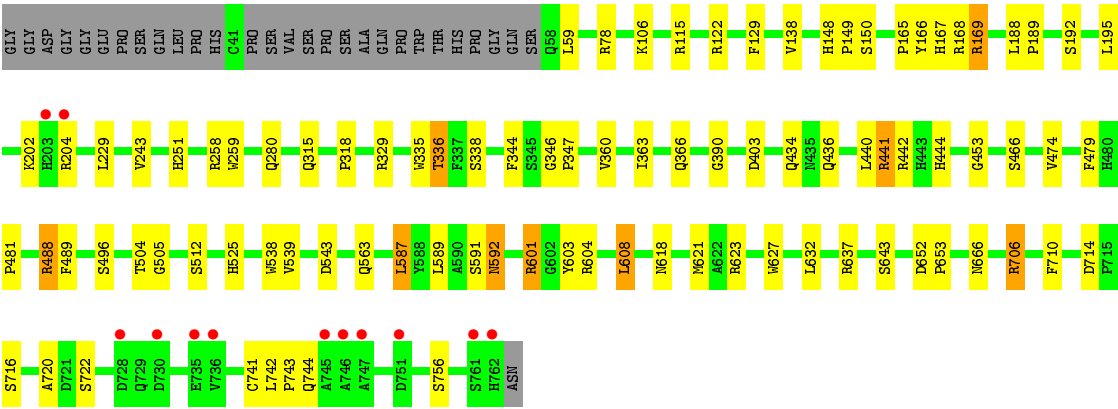


• Molecule 1: MEMBRANE COPPER AMINE OXIDASE

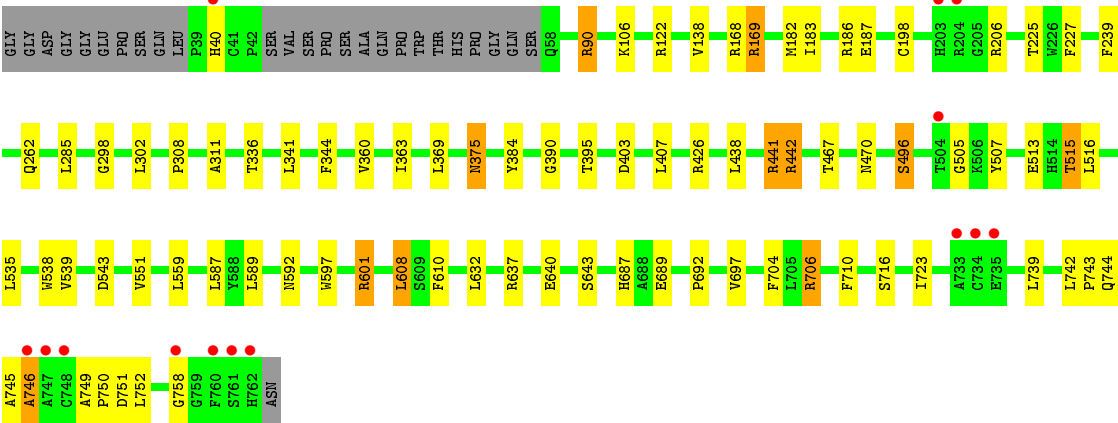
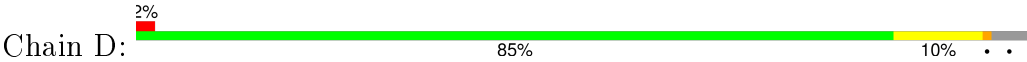


• Molecule 1: MEMBRANE COPPER AMINE OXIDASE





● Molecule 1: MEMBRANE COPPER AMINE OXIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	130.24Å 130.24Å 221.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 19.97 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-2.50) 98.2 (19.97-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.204 , 0.252 0.204 , 0.206	Depositor DCC
R_{free} test set	6351 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.645	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 12.4	EDS
Estimated twinning fraction	0.448 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 126403 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23470	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, CL, CA, NDG, TPQ, FUL, CU, MAN

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	0.52	0/5724	0.64	1/7804 (0.0%)
1	B	0.51	0/5751	0.62	1/7842 (0.0%)
1	C	0.51	0/5724	0.62	1/7804 (0.0%)
1	D	0.51	0/5751	0.63	0/7842
All	All	0.51	0/22950	0.63	3/31292 (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	D	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	341	LEU	CA-CB-CG	5.25	127.38	115.30
1	B	589	LEU	CA-CB-CG	5.05	126.93	115.30
1	C	592	ASN	N-CA-CB	-5.00	101.60	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	198	CYS	Peptide

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	5563	0	5311	69	0
1	B	5587	0	5333	49	0
1	C	5563	0	5311	49	0
1	D	5587	0	5332	57	0
2	A	39	0	34	0	0
2	B	39	0	34	1	0
2	C	39	0	34	0	0
3	A	38	0	34	0	0
3	C	38	0	34	1	0
4	A	28	0	25	0	0
4	B	28	0	25	0	0
4	C	28	0	25	0	0
4	D	28	0	25	0	0
5	A	14	0	13	1	0
5	B	42	0	39	1	0
5	C	14	0	13	0	0
5	D	42	0	39	1	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
6	C	2	0	0	0	0
6	D	2	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	A	2	0	0	0	0
8	B	2	0	0	0	0
8	C	2	0	0	0	0
8	D	2	0	0	0	0
9	B	60	0	52	2	0
9	D	60	0	52	0	0
10	D	50	0	43	0	0
11	A	141	0	0	9	0
11	B	139	0	0	4	0
11	C	141	0	0	5	0
11	D	142	0	0	5	0
All	All	23470	0	21808	209	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 5.

All (209) 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:D:360:VAL:HG11	1:D:363:ILE:HG13	1.38	1.04
1:A:360:VAL:HG11	1:A:363:ILE:HG13	1.56	0.85
1:D:426:ARG:HE	1:D:758:GLY:HA3	1.49	0.78
1:A:538:TRP:CZ3	1:A:592:ASN:HB2	2.19	0.77
1:A:488:ARG:HH12	1:A:608:LEU:CD1	1.98	0.77
1:C:488:ARG:HH22	1:C:608:LEU:HD13	1.48	0.76
1:C:360:VAL:HG11	1:C:363:ILE:HG13	1.69	0.74
1:A:488:ARG:HH12	1:A:608:LEU:HD13	1.52	0.74
1:A:166:TYR:O	1:A:169:ARG:HG3	1.88	0.74
1:B:106:LYS:HB2	1:B:637:ARG:NH2	2.04	0.72
1:A:403:ASP:OD1	1:B:442:ARG:HG3	1.89	0.72
1:D:587:LEU:HD22	1:D:632:LEU:HD21	1.70	0.71
1:C:166:TYR:O	1:C:169:ARG:HG3	1.92	0.70
1:A:453:GLY:HA3	1:B:302:LEU:HD13	1.77	0.67
1:A:706:ARG:CZ	11:A:2137:HOH:O	2.44	0.66
1:A:706:ARG:NE	11:A:2137:HOH:O	2.28	0.66
1:C:403:ASP:OD1	1:D:442:ARG:HG3	1.95	0.66
1:A:587:LEU:HD22	1:A:632:LEU:HD21	1.78	0.66
1:A:315:GLN:HE22	1:A:434:GLN:HA	1.61	0.65
1:A:539:VAL:HG22	1:A:589:LEU:HD22	1.77	0.65
1:D:407:LEU:HD21	1:D:752:LEU:HD23	1.79	0.64
1:C:538:TRP:CZ3	1:C:592:ASN:HB2	2.34	0.63
1:B:587:LEU:HD22	1:B:632:LEU:HD21	1.80	0.63
1:B:285:LEU:O	1:B:285:LEU:HG	1.99	0.63
1:B:539:VAL:HG22	1:B:589:LEU:HD22	1.81	0.62
5:D:1774:NAG:O4	11:D:2142:HOH:O	2.13	0.62
1:D:106:LYS:HB2	1:D:637:ARG:NH2	2.14	0.62
1:C:488:ARG:HH12	1:C:608:LEU:CD1	2.11	0.62
1:D:344:PHE:HA	1:D:390:GLY:HA2	1.80	0.62
1:B:468:LEU:HB2	1:B:471:TPQ:O5	2.00	0.62
1:A:441:ARG:NH2	11:A:2084:HOH:O	2.34	0.61
1:C:329:ARG:HH21	1:C:336:THR:HG23	1.66	0.60
9:B:1765:NAG:H5	9:B:1766:FUL:H61	1.84	0.60
1:A:640:GLU:CD	1:A:640:GLU:H	2.05	0.60
1:A:315:GLN:NE2	1:A:434:GLN:HA	2.16	0.59
1:A:98:SER:HB2	1:A:417:LEU:HD23	1.84	0.59
1:B:723:ILE:HD13	1:B:739:LEU:HD22	1.84	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:PHE:CZ	1:A:169:ARG:HB2	2.37	0.59
1:C:587:LEU:HD22	1:C:632:LEU:HD21	1.83	0.59
1:A:344:PHE:HA	1:A:390:GLY:HA2	1.85	0.59
1:D:169:ARG:HD2	11:D:2018:HOH:O	2.03	0.59
1:A:302:LEU:HD13	1:B:453:GLY:HA3	1.85	0.58
1:A:608:LEU:HD23	1:B:610:PHE:HZ	1.68	0.58
1:A:706:ARG:NH2	11:A:2137:HOH:O	2.37	0.58
9:B:1765:NAG:O3	5:B:1773:NAG:H2	2.04	0.58
1:D:742:LEU:O	1:D:746:ALA:HB3	2.05	0.57
1:C:315:GLN:HE22	1:C:434:GLN:HA	1.70	0.57
1:A:591:SER:OG	1:A:592:ASN:N	2.32	0.57
1:C:442:ARG:HG3	1:D:403:ASP:OD1	2.04	0.57
1:A:366:GLN:OE1	1:A:643:SER:HA	2.06	0.56
1:C:714:ASP:HB2	1:D:689:GLU:O	2.06	0.56
1:D:539:VAL:HG22	1:D:589:LEU:HD22	1.88	0.56
1:C:436:GLN:OE1	1:C:436:GLN:HA	2.05	0.55
1:A:742:LEU:N	1:A:743:PRO:HD2	2.22	0.55
1:D:587:LEU:CD2	1:D:632:LEU:HD21	2.35	0.55
1:C:539:VAL:HG22	1:C:589:LEU:HD22	1.89	0.54
1:D:360:VAL:HG11	1:D:363:ILE:CG1	2.25	0.54
1:D:587:LEU:HD22	1:D:632:LEU:CD2	2.35	0.54
1:B:440:LEU:HD23	1:B:455:LEU:HD23	1.88	0.54
1:C:453:GLY:HA3	1:D:302:LEU:HD13	1.89	0.54
1:B:538:TRP:CZ3	1:B:592:ASN:HB2	2.42	0.54
1:C:188:LEU:HD22	1:C:195:LEU:HD11	1.90	0.54
1:D:608:LEU:HD11	1:D:704:PHE:CD2	2.42	0.54
1:D:723:ILE:HD13	1:D:739:LEU:HD22	1.89	0.54
1:A:488:ARG:HH22	1:A:608:LEU:HD13	1.73	0.53
1:B:513:GLU:O	1:B:515:THR:HG22	2.07	0.53
1:D:298:GLY:O	1:D:692:PRO:HB3	2.09	0.53
1:B:90:ARG:HG3	1:B:91:PRO:HD2	1.91	0.53
1:B:392:GLY:O	1:B:395:THR:HB	2.09	0.53
1:C:706:ARG:NE	11:C:2128:HOH:O	2.41	0.53
1:D:285:LEU:HG	1:D:285:LEU:O	2.09	0.53
1:A:315:GLN:NE2	1:A:435:ASN:H	2.07	0.52
1:C:441:ARG:NH2	11:C:2072:HOH:O	2.41	0.52
1:B:168:ARG:HD3	11:B:2124:HOH:O	2.09	0.52
1:B:643:SER:HB3	11:B:2053:HOH:O	2.09	0.52
1:C:440:LEU:HD22	1:C:481:PRO:HG2	1.91	0.51
1:B:400:ARG:HD3	1:B:406:TYR:CE2	2.45	0.51
1:A:441:ARG:HD3	1:A:716:SER:OG	2.10	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:SER:O	1:A:380:MET:HB2	2.11	0.51
1:B:329:ARG:HH21	1:B:336:THR:HG23	1.76	0.51
1:B:413:TRP:HE3	1:B:425:ILE:HD12	1.76	0.51
1:D:441:ARG:HD3	1:D:716:SER:OG	2.11	0.50
1:D:687:HIS:HD1	1:D:689:GLU:HG2	1.76	0.50
1:A:78:ARG:HB3	1:A:78:ARG:HH11	1.75	0.50
1:C:344:PHE:HA	1:C:390:GLY:HA2	1.92	0.50
1:D:608:LEU:HD11	1:D:704:PHE:CE2	2.47	0.50
1:C:621:MET:HB2	1:C:653:PRO:HB2	1.93	0.49
1:A:188:LEU:N	1:A:189:PRO:HD2	2.26	0.49
1:C:608:LEU:HD23	1:D:610:PHE:HZ	1.77	0.49
1:A:720:ALA:O	1:B:308:PRO:HA	2.13	0.49
1:A:85:ASP:CG	1:A:420:GLN:HG2	2.33	0.49
1:D:640:GLU:CD	1:D:640:GLU:H	2.15	0.49
1:C:129:PHE:CZ	1:C:169:ARG:HB2	2.48	0.48
3:C:1764:NDG:H6C1	3:C:1765:NAG:C1	2.44	0.48
1:C:563:GLN:HG2	1:D:507:TYR:CE1	2.48	0.48
1:B:601:ARG:HB3	1:B:710:PHE:HA	1.96	0.48
1:B:59:LEU:HD11	1:B:120:PRO:HD2	1.95	0.48
1:D:106:LYS:HE2	1:D:535:LEU:HD21	1.96	0.48
1:C:441:ARG:HD3	1:C:716:SER:OG	2.14	0.48
1:A:444:HIS:O	1:B:467:THR:HG21	2.13	0.48
1:C:318:PRO:HG2	1:D:311:ALA:HB3	1.96	0.47
1:A:714:ASP:HB2	1:B:689:GLU:O	2.12	0.47
1:C:601:ARG:HB3	1:C:710:PHE:HA	1.96	0.47
1:D:496:SER:HB3	11:D:2084:HOH:O	2.15	0.47
1:D:637:ARG:HG3	1:D:637:ARG:HH11	1.79	0.47
1:A:743:PRO:HA	1:B:752:LEU:HD13	1.96	0.47
1:A:98:SER:HB2	1:A:417:LEU:CD2	2.45	0.47
1:B:181:GLN:NE2	1:B:185:ASN:OD1	2.47	0.47
1:C:488:ARG:HH12	1:C:608:LEU:HD12	1.80	0.47
1:B:298:GLY:O	1:B:692:PRO:HB3	2.15	0.47
1:D:168:ARG:HD3	11:D:2118:HOH:O	2.15	0.47
1:D:687:HIS:ND1	1:D:689:GLU:HG2	2.29	0.46
1:C:504:THR:HG22	1:C:505:GLY:H	1.80	0.46
1:A:106:LYS:HB2	1:A:637:ARG:NH1	2.31	0.46
1:B:203:HIS:H	1:B:203:HIS:CD2	2.33	0.46
1:C:720:ALA:O	1:D:308:PRO:HA	2.15	0.46
1:A:442:ARG:HG2	1:A:443:HIS:N	2.25	0.46
1:A:488:ARG:NH1	1:A:608:LEU:HD13	2.24	0.45
1:B:329:ARG:HH21	1:B:336:THR:CG2	2.29	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:513:GLU:O	1:D:515:THR:HG23	2.16	0.45
1:D:262:GLN:HG2	11:D:2034:HOH:O	2.16	0.45
1:C:488:ARG:NH2	1:C:608:LEU:HD13	2.25	0.45
1:A:440:LEU:HD22	1:A:481:PRO:HG2	1.98	0.45
1:D:551:VAL:HG22	1:D:559:LEU:HD23	1.97	0.45
1:C:366:GLN:OE1	1:C:643:SER:HA	2.17	0.45
1:D:239:PHE:CD2	1:D:470:ASN:HB3	2.52	0.45
1:D:375:ASN:HD22	1:D:375:ASN:HA	1.55	0.45
1:A:593:HIS:O	1:A:601:ARG:HG2	2.17	0.45
1:D:183:ILE:HA	1:D:187:GLU:HB2	1.99	0.45
1:B:551:VAL:HG22	1:B:559:LEU:HD23	1.99	0.45
2:B:1763:NAG:H4	2:B:1769:BMA:O2	2.17	0.44
1:D:704:PHE:HB3	1:D:706:ARG:HD3	1.99	0.44
1:C:706:ARG:CZ	11:C:2128:HOH:O	2.65	0.44
1:B:224:ALA:HA	1:B:249:VAL:O	2.17	0.44
1:C:525:HIS:HB2	1:C:627:TRP:CE3	2.53	0.44
1:D:601:ARG:CB	1:D:710:PHE:HA	2.47	0.44
1:D:538:TRP:CZ3	1:D:592:ASN:HB2	2.53	0.44
1:A:122:ARG:HB3	1:A:146:LEU:HB2	2.00	0.43
1:A:59:LEU:HD11	1:A:120:PRO:HD2	2.00	0.43
1:D:745:ALA:O	1:D:746:ALA:HB2	2.18	0.43
1:D:225:THR:HB	1:D:227:PHE:CE1	2.53	0.43
1:B:267:GLN:NE2	11:B:2035:HOH:O	2.42	0.43
1:A:335:TRP:CE2	1:A:479:PHE:HB3	2.54	0.43
1:A:393:LYS:HG3	1:A:417:LEU:HD13	2.00	0.43
1:C:168:ARG:HG2	1:C:652:ASP:HB2	2.00	0.43
1:A:739:LEU:HD21	1:B:401:GLY:HA3	2.01	0.43
1:A:604:ARG:HD3	11:A:2136:HOH:O	2.18	0.43
1:B:189:PRO:O	1:B:192:SER:HB2	2.19	0.43
1:A:90:ARG:NH2	11:A:2007:HOH:O	2.49	0.43
1:B:350:PHE:CE2	1:B:362:GLU:HG3	2.53	0.43
1:D:369:LEU:HD12	1:D:384:TYR:O	2.19	0.43
1:A:114:ASP:OD2	1:A:358:ARG:NH2	2.48	0.43
1:A:400:ARG:HH21	1:A:400:ARG:HB2	1.84	0.42
1:C:165:PRO:HB2	1:C:167:HIS:CE1	2.53	0.42
1:A:512:SER:HA	1:B:597:TRP:CZ2	2.55	0.42
1:B:608:LEU:HD11	1:B:704:PHE:CD2	2.54	0.42
1:B:584:PRO:HG2	1:B:587:LEU:HD12	2.01	0.42
1:A:61:ALA:O	1:A:122:ARG:NH2	2.53	0.42
1:B:262:GLN:HG2	11:B:2033:HOH:O	2.18	0.42
1:A:191:ALA:HA	1:A:278:GLU:HB2	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:GLY:O	1:A:692:PRO:HB3	2.19	0.42
1:A:241:HIS:NE2	1:A:299:SER:O	2.52	0.42
1:A:525:HIS:HB2	1:A:627:TRP:CE3	2.53	0.42
1:C:335:TRP:CE2	1:C:479:PHE:HB3	2.54	0.42
1:A:495:ILE:HG12	11:A:2094:HOH:O	2.19	0.42
1:D:40:HIS:ND1	1:D:751:ASP:HB3	2.35	0.42
1:C:742:LEU:N	1:C:743:PRO:HD2	2.35	0.42
1:D:341:LEU:HD21	1:D:395:THR:OG1	2.19	0.42
1:B:360:VAL:HG22	1:B:530:LEU:HD23	2.02	0.42
1:A:360:VAL:HG11	1:A:363:ILE:CG1	2.39	0.42
1:C:706:ARG:NH2	11:C:2128:HOH:O	2.52	0.42
1:A:107:ALA:HB3	11:A:2013:HOH:O	2.19	0.42
1:D:90:ARG:HD3	1:D:90:ARG:HA	1.61	0.42
1:A:181:GLN:O	1:A:185:ASN:HB2	2.19	0.42
1:C:189:PRO:O	1:C:192:SER:HB2	2.20	0.42
1:C:512:SER:HA	1:D:597:TRP:CZ2	2.55	0.42
1:C:346:GLY:HA3	1:C:347:PRO:HD2	1.90	0.42
1:A:78:ARG:HB3	1:A:78:ARG:NH1	2.34	0.41
1:A:665:ASN:OD1	1:A:665:ASN:N	2.54	0.41
1:A:198:CYS:HB3	1:A:288:VAL:HG11	2.01	0.41
1:A:351:ASP:OD2	1:A:358:ARG:HD3	2.20	0.41
1:A:599:HIS:HA	1:A:600:PRO:HD3	1.97	0.41
1:B:370:ALA:HA	1:B:521:THR:O	2.20	0.41
1:D:535:LEU:HA	1:D:535:LEU:HD23	1.80	0.41
1:B:687:HIS:HD1	1:B:689:GLU:HG2	1.86	0.41
1:A:165:PRO:HB2	1:A:167:HIS:CE1	2.55	0.41
1:D:742:LEU:HB3	1:D:743:PRO:HD3	2.01	0.41
1:D:687:HIS:CE1	1:D:689:GLU:HG2	2.55	0.41
1:C:106:LYS:HB2	1:C:637:ARG:NH1	2.36	0.41
1:C:474:VAL:O	1:C:489:PHE:HA	2.21	0.41
1:B:138:VAL:O	1:B:164:LEU:HB2	2.20	0.41
1:C:444:HIS:O	1:D:467:THR:HG21	2.21	0.41
1:C:251:HIS:HA	1:C:259:TRP:CD1	2.56	0.41
1:B:737:ASN:HA	1:B:738:PRO:HD2	1.81	0.40
5:A:1768:NAG:H61	11:A:2141:HOH:O	2.20	0.40
1:D:749:ALA:HA	1:D:750:PRO:HD2	1.83	0.40
1:C:148:HIS:HA	1:C:149:PRO:HD3	1.80	0.40
1:C:743:PRO:HA	1:D:752:LEU:HD12	2.02	0.40
1:D:182:MET:HG3	1:D:186:ARG:NH2	2.35	0.40
1:A:663:PHE:N	1:A:663:PHE:CD1	2.90	0.40
1:C:604:ARG:HD3	11:C:2109:HOH:O	2.20	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:591:SER:OG	1:C:592:ASN:N	2.55	0.40
1:D:601:ARG:HB2	1:D:710:PHE:HA	2.02	0.40
1:B:624:GLY:HA2	1:B:656:PRO:HB3	2.03	0.40
1:B:90:ARG:HD3	1:B:90:ARG:HA	1.90	0.40
1:A:189:PRO:O	1:A:192:SER:HB2	2.21	0.40
1:A:525:HIS:CE1	1:A:677:TRP:HB3	2.57	0.40
1:B:376:SER:O	1:B:380:MET:HB2	2.22	0.40
1:B:663:PHE:CD1	1:B:663:PHE:N	2.89	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	702/735 (96%)	671 (96%)	30 (4%)	1 (0%)	56	78
1	B	704/735 (96%)	675 (96%)	27 (4%)	2 (0%)	46	68
1	C	702/735 (96%)	668 (95%)	32 (5%)	2 (0%)	46	68
1	D	704/735 (96%)	675 (96%)	26 (4%)	3 (0%)	39	61
All	All	2812/2940 (96%)	2689 (96%)	115 (4%)	8 (0%)	46	68

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	746	ALA
1	C	744	GLN
1	B	59	LEU
1	C	204	ARG
1	B	205	GLY
1	D	744	GLN
1	A	738	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	505	GLY

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	585/609 (96%)	555 (95%)	30 (5%)	29	52
1	B	588/609 (97%)	568 (97%)	20 (3%)	44	72
1	C	585/609 (96%)	555 (95%)	30 (5%)	29	52
1	D	588/609 (97%)	569 (97%)	19 (3%)	46	74
All	All	2346/2436 (96%)	2247 (96%)	99 (4%)	36	62

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	LEU
1	A	115	ARG
1	A	122	ARG
1	A	138	VAL
1	A	150	SER
1	A	169	ARG
1	A	202	LYS
1	A	229	LEU
1	A	258	ARG
1	A	280	GLN
1	A	336	THR
1	A	395	THR
1	A	441	ARG
1	A	442	ARG
1	A	464	SER
1	A	466	SER
1	A	489	PHE
1	A	496	SER
1	A	549	MET
1	A	587	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	601	ARG
1	A	603	TYR
1	A	608	LEU
1	A	618	ASN
1	A	637	ARG
1	A	640	GLU
1	A	665	ASN
1	A	666	ASN
1	A	722	SER
1	A	741	CYS
1	B	40	HIS
1	B	76	THR
1	B	90	ARG
1	B	138	VAL
1	B	229	LEU
1	B	336	THR
1	B	375	ASN
1	B	395	THR
1	B	441	ARG
1	B	515	THR
1	B	556	GLU
1	B	587	LEU
1	B	601	ARG
1	B	608	LEU
1	B	640	GLU
1	B	665	ASN
1	B	693	ASN
1	B	697	VAL
1	B	706	ARG
1	B	722	SER
1	C	59	LEU
1	C	78	ARG
1	C	115	ARG
1	C	122	ARG
1	C	138	VAL
1	C	150	SER
1	C	169	ARG
1	C	202	LYS
1	C	229	LEU
1	C	243	VAL
1	C	258	ARG
1	C	280	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	336	THR
1	C	338	SER
1	C	441	ARG
1	C	466	SER
1	C	488	ARG
1	C	496	SER
1	C	543	ASP
1	C	587	LEU
1	C	601	ARG
1	C	603	TYR
1	C	608	LEU
1	C	618	ASN
1	C	623	ARG
1	C	666	ASN
1	C	706	ARG
1	C	722	SER
1	C	741	CYS
1	C	756	SER
1	D	90	ARG
1	D	122	ARG
1	D	138	VAL
1	D	169	ARG
1	D	206	ARG
1	D	336	THR
1	D	375	ASN
1	D	438	LEU
1	D	441	ARG
1	D	442	ARG
1	D	496	SER
1	D	515	THR
1	D	516	LEU
1	D	543	ASP
1	D	601	ARG
1	D	608	LEU
1	D	643	SER
1	D	697	VAL
1	D	706	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	242	HIS
1	A	262	GLN
1	A	319	GLN
1	A	563	GLN
1	A	666	ASN
1	B	40	HIS
1	B	181	GLN
1	B	203	HIS
1	B	618	ASN
1	C	185	ASN
1	D	375	ASN
1	D	420	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	TPQ	A	471	1	13,14,15	2.08	3 (23%)	15,19,21	2.80	6 (40%)
1	TPQ	B	471	1	13,14,15	2.13	3 (23%)	15,19,21	2.74	5 (33%)
1	TPQ	C	471	1	13,14,15	2.10	3 (23%)	15,19,21	2.55	4 (26%)
1	TPQ	D	471	1	13,14,15	2.11	3 (23%)	15,19,21	2.56	5 (33%)

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	TPQ	A	471	1	-	0/4/22/24	0/1/1/1
1	TPQ	B	471	1	-	0/4/22/24	0/1/1/1
1	TPQ	C	471	1	-	0/4/22/24	0/1/1/1
1	TPQ	D	471	1	-	0/4/22/24	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	471	TPQ	C3-C4	2.07	1.39	1.35
1	C	471	TPQ	C3-C4	2.36	1.39	1.35
1	A	471	TPQ	C3-C4	2.49	1.39	1.35
1	B	471	TPQ	C3-C4	2.64	1.40	1.35
1	B	471	TPQ	O2-C2	4.28	1.36	1.24
1	D	471	TPQ	O2-C2	4.28	1.36	1.24
1	A	471	TPQ	O2-C2	4.29	1.36	1.24
1	C	471	TPQ	O2-C2	4.35	1.36	1.24
1	A	471	TPQ	O5-C5	4.43	1.36	1.24
1	B	471	TPQ	O5-C5	4.61	1.37	1.24
1	C	471	TPQ	O5-C5	4.68	1.37	1.24
1	D	471	TPQ	O5-C5	4.68	1.37	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	471	TPQ	C1-C6-C5	-7.95	118.13	122.97
1	B	471	TPQ	C1-C6-C5	-7.49	118.41	122.97
1	C	471	TPQ	C1-C6-C5	-7.33	118.50	122.97
1	D	471	TPQ	C1-C6-C5	-6.45	119.04	122.97
1	B	471	TPQ	O2-C2-C1	-4.22	117.16	120.85
1	D	471	TPQ	O2-C2-C3	-2.40	116.48	121.89
1	A	471	TPQ	O2-C2-C1	-2.36	118.79	120.85
1	C	471	TPQ	O2-C2-C1	-2.15	118.97	120.85
1	A	471	TPQ	C4-C3-C2	-2.07	118.14	120.77
1	B	471	TPQ	C6-C5-C4	2.30	121.28	117.34
1	A	471	TPQ	C6-C1-C2	2.37	120.11	118.44
1	A	471	TPQ	C6-C5-C4	2.38	121.43	117.34
1	D	471	TPQ	C6-C5-C4	2.48	121.59	117.34
1	D	471	TPQ	CA-CB-C1	2.77	119.17	113.63
1	C	471	TPQ	C6-C1-C2	3.39	120.83	118.44
1	B	471	TPQ	C6-C1-C2	3.52	120.92	118.44
1	B	471	TPQ	C3-C2-C1	3.60	120.99	118.30
1	C	471	TPQ	C3-C2-C1	3.80	121.14	118.30
1	A	471	TPQ	C3-C2-C1	4.85	121.93	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	471	TPQ	C3-C2-C1	4.91	121.98	118.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	471	TPQ	1	0

5.5 Carbohydrates ⓘ

37 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	1762	1,2	14,14,15	0.59	0	15,19,21	1.43	3 (20%)
2	NAG	A	1763	2	14,14,15	0.63	0	15,19,21	1.27	1 (6%)
3	NDG	A	1764	1,3	14,14,15	0.65	0	15,19,21	1.59	3 (20%)
3	NAG	A	1765	3	14,14,15	0.63	0	15,19,21	1.63	3 (20%)
3	FUL	A	1766	3	10,10,11	0.68	0	14,14,16	2.12	2 (14%)
4	NAG	A	1767	1,4	14,14,15	0.55	0	15,19,21	1.03	1 (6%)
2	BMA	A	1769	2	11,11,12	0.48	0	14,15,17	1.31	1 (7%)
4	NAG	A	1770	4	14,14,15	0.60	0	15,19,21	1.37	1 (6%)
2	NAG	B	1762	1,2	14,14,15	0.64	0	15,19,21	1.33	2 (13%)
2	NAG	B	1763	2	14,14,15	0.48	0	15,19,21	1.13	3 (20%)
9	NDG	B	1764	1,9	14,14,15	0.57	0	15,19,21	1.53	1 (6%)
9	NAG	B	1765	9	14,14,15	0.51	0	15,19,21	1.07	1 (6%)
9	FUL	B	1766	9	10,10,11	0.68	0	14,14,16	1.18	1 (7%)
4	NAG	B	1767	1,4	14,14,15	0.62	0	15,19,21	0.98	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BMA	B	1769	2	11,11,12	0.75	0	14,15,17	0.91	1 (7%)
4	NAG	B	1771	4	14,14,15	0.58	0	15,19,21	1.70	2 (13%)
9	BMA	B	1772	9	11,11,12	0.71	0	14,15,17	2.08	3 (21%)
9	MAN	B	1774	9	11,11,12	0.55	0	14,15,17	1.26	1 (7%)
2	NAG	C	1762	1,2	14,14,15	0.55	0	15,19,21	0.97	1 (6%)
2	NAG	C	1763	2	14,14,15	0.52	0	15,19,21	1.04	1 (6%)
3	NDG	C	1764	1,3	14,14,15	0.61	0	15,19,21	1.82	2 (13%)
3	NAG	C	1765	3	14,14,15	0.53	0	15,19,21	1.22	2 (13%)
3	FUL	C	1766	3	10,10,11	0.79	0	14,14,16	1.83	4 (28%)
4	NAG	C	1767	1,4	14,14,15	0.52	0	15,19,21	1.43	2 (13%)
2	BMA	C	1769	2	11,11,12	0.60	0	14,15,17	0.91	1 (7%)
4	NAG	C	1770	4	14,14,15	0.61	0	15,19,21	1.77	1 (6%)
10	NAG	D	1762	1,10	14,14,15	0.54	0	15,19,21	0.99	0
10	NAG	D	1763	10	14,14,15	0.54	0	15,19,21	2.00	6 (40%)
9	NDG	D	1764	1,9	14,14,15	0.49	0	15,19,21	1.98	3 (20%)
9	NAG	D	1765	9	14,14,15	0.48	0	15,19,21	1.96	2 (13%)
9	FUL	D	1766	9	10,10,11	0.75	0	14,14,16	1.79	3 (21%)
4	NAG	D	1767	1,4	14,14,15	0.56	0	15,19,21	0.90	1 (6%)
10	BMA	D	1769	10	11,11,12	0.56	0	14,15,17	1.21	1 (7%)
4	NAG	D	1771	4	14,14,15	0.61	0	15,19,21	1.19	1 (6%)
9	BMA	D	1772	9	11,11,12	0.92	1 (9%)	14,15,17	2.32	3 (21%)
10	MAN	D	1773	10	11,11,12	0.62	0	14,15,17	1.51	2 (14%)
9	MAN	D	1775	9	11,11,12	0.57	0	14,15,17	2.79	3 (21%)

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	1762	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1763	2	-	0/6/23/26	0/1/1/1
3	NDG	A	1764	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1765	3	-	0/6/23/26	0/1/1/1
3	FUL	A	1766	3	-	0/0/17/20	0/1/1/1
4	NAG	A	1767	1,4	-	0/6/23/26	0/1/1/1
2	BMA	A	1769	2	-	0/2/19/22	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1770	4	-	0/6/23/26	0/1/1/1
2	NAG	B	1762	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	1763	2	-	0/6/23/26	0/1/1/1
9	NDG	B	1764	1,9	-	0/6/23/26	0/1/1/1
9	NAG	B	1765	9	-	0/6/23/26	0/1/1/1
9	FUL	B	1766	9	-	0/0/17/20	0/1/1/1
4	NAG	B	1767	1,4	-	0/6/23/26	0/1/1/1
2	BMA	B	1769	2	-	0/2/19/22	0/1/1/1
4	NAG	B	1771	4	-	0/6/23/26	0/1/1/1
9	BMA	B	1772	9	-	0/2/19/22	0/1/1/1
9	MAN	B	1774	9	-	0/2/19/22	0/1/1/1
2	NAG	C	1762	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	1763	2	-	0/6/23/26	0/1/1/1
3	NDG	C	1764	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	1765	3	-	0/6/23/26	0/1/1/1
3	FUL	C	1766	3	-	0/0/17/20	0/1/1/1
4	NAG	C	1767	1,4	-	0/6/23/26	0/1/1/1
2	BMA	C	1769	2	-	0/2/19/22	0/1/1/1
4	NAG	C	1770	4	-	0/6/23/26	0/1/1/1
10	NAG	D	1762	1,10	-	0/6/23/26	0/1/1/1
10	NAG	D	1763	10	-	0/6/23/26	0/1/1/1
9	NDG	D	1764	1,9	-	0/6/23/26	0/1/1/1
9	NAG	D	1765	9	-	0/6/23/26	0/1/1/1
9	FUL	D	1766	9	-	0/0/17/20	0/1/1/1
4	NAG	D	1767	1,4	-	0/6/23/26	0/1/1/1
10	BMA	D	1769	10	-	0/2/19/22	0/1/1/1
4	NAG	D	1771	4	-	0/6/23/26	0/1/1/1
9	BMA	D	1772	9	-	0/2/19/22	0/1/1/1
10	MAN	D	1773	10	-	0/2/19/22	0/1/1/1
9	MAN	D	1775	9	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	D	1772	BMA	C2-C3	2.18	1.55	1.52

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1763	NAG	C2-N2-C7	-3.98	117.92	123.04
9	D	1764	NDG	O7-C7-C8	-2.60	117.29	122.06
3	C	1764	NDG	O6-C6-C5	-2.50	103.08	111.33

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1762	NAG	O4-C4-C3	-2.46	104.80	110.34
10	D	1763	NAG	O7-C7-C8	-2.45	117.57	122.06
10	D	1763	NAG	C6-C5-C4	-2.42	107.06	113.02
9	D	1775	MAN	C2-C3-C4	-2.41	106.95	111.04
2	A	1762	NAG	O4-C4-C3	-2.36	105.02	110.34
2	B	1762	NAG	O4-C4-C3	-2.20	105.38	110.34
2	C	1763	NAG	C4-C3-C2	-2.20	107.81	111.23
4	A	1767	NAG	O4-C4-C3	-2.20	105.39	110.34
9	D	1766	FUL	C2-C3-C4	-2.18	107.34	111.04
4	B	1771	NAG	C4-C3-C2	-2.15	107.88	111.23
3	A	1764	NDG	O-C5-C6	-2.10	102.79	107.35
2	A	1762	NAG	C2-N2-C7	-2.07	120.39	123.04
9	D	1764	NDG	O6-C6-C5	-2.05	104.57	111.33
2	B	1763	NAG	O4-C4-C3	-2.04	105.75	110.34
3	A	1764	NDG	O6-C6-C5	-2.03	104.63	111.33
2	B	1763	NAG	C1-O5-C5	2.02	114.81	112.25
2	B	1769	BMA	O5-C5-C6	2.09	111.87	107.35
4	D	1767	NAG	C1-O5-C5	2.09	114.90	112.25
2	C	1769	BMA	C1-O5-C5	2.12	114.94	112.25
3	C	1765	NAG	C4-C3-C2	2.16	114.59	111.23
3	C	1766	FUL	O5-C5-C6	2.25	109.85	106.13
3	C	1766	FUL	O5-C5-C4	2.25	113.44	109.53
9	B	1772	BMA	C1-O5-C5	2.29	115.15	112.25
2	B	1763	NAG	C4-C3-C2	2.31	114.81	111.23
10	D	1763	NAG	C2-N2-C7	2.37	126.09	123.04
9	B	1766	FUL	O5-C5-C6	2.41	110.11	106.13
4	C	1767	NAG	C1-O5-C5	2.43	115.33	112.25
9	D	1772	BMA	C3-C4-C5	2.45	114.46	110.20
3	C	1765	NAG	O5-C5-C6	2.53	112.83	107.35
9	B	1765	NAG	C1-O5-C5	2.61	115.56	112.25
3	A	1765	NAG	O5-C5-C6	2.66	113.10	107.35
10	D	1763	NAG	C3-C4-C5	2.67	114.84	110.20
9	D	1765	NAG	C2-N2-C7	2.76	126.59	123.04
10	D	1763	NAG	C8-C7-N2	2.76	121.39	116.11
2	B	1762	NAG	C1-O5-C5	2.78	115.77	112.25
10	D	1773	MAN	C1-C2-C3	2.95	113.03	109.54
3	A	1765	NAG	C4-C3-C2	3.03	115.94	111.23
4	B	1767	NAG	C4-C3-C2	3.06	115.99	111.23
9	B	1772	BMA	C2-C3-C4	3.11	116.33	111.04
4	D	1771	NAG	C1-O5-C5	3.19	116.30	112.25
3	A	1766	FUL	C1-C2-C3	3.37	113.53	109.54
9	D	1772	BMA	C2-C3-C4	3.38	116.78	111.04

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1762	NAG	C1-O5-C5	3.39	116.56	112.25
4	C	1767	NAG	C2-N2-C7	3.43	127.45	123.04
3	C	1766	FUL	C1-C2-C3	3.54	113.73	109.54
9	D	1766	FUL	O5-C5-C6	3.67	112.20	106.13
9	D	1766	FUL	C1-O5-C5	3.77	118.19	112.38
9	D	1775	MAN	O5-C1-C2	3.83	117.06	110.86
10	D	1769	BMA	C1-O5-C5	3.86	117.14	112.25
10	D	1773	MAN	C1-O5-C5	3.87	117.16	112.25
4	A	1770	NAG	C1-O5-C5	3.88	117.17	112.25
3	A	1765	NAG	C1-O5-C5	3.96	117.28	112.25
9	B	1774	MAN	C1-O5-C5	3.96	117.28	112.25
3	C	1766	FUL	C1-O5-C5	3.97	118.52	112.38
2	A	1769	BMA	C1-O5-C5	4.16	117.52	112.25
10	D	1763	NAG	C1-O5-C5	4.18	117.55	112.25
9	B	1764	NDG	C1-O-C5	4.50	117.95	112.25
3	A	1764	NDG	C1-O-C5	4.51	117.97	112.25
4	B	1771	NAG	C1-O5-C5	4.93	118.51	112.25
3	C	1764	NDG	C1-O-C5	5.31	118.98	112.25
4	C	1770	NAG	C1-O5-C5	5.81	119.62	112.25
9	B	1772	BMA	C1-C2-C3	5.84	116.45	109.54
3	A	1766	FUL	C1-O5-C5	5.99	121.63	112.38
9	D	1764	NDG	C1-O-C5	6.12	120.02	112.25
9	D	1765	NAG	C1-O5-C5	6.28	120.22	112.25
9	D	1772	BMA	C1-C2-C3	7.15	118.00	109.54
9	D	1775	MAN	C1-O5-C5	8.81	123.42	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1763	NAG	1	0
9	B	1765	NAG	2	0
9	B	1766	FUL	1	0
2	B	1769	BMA	1	0
3	C	1764	NDG	1	0
3	C	1765	NAG	1	0

5.6 Ligand geometry

Of 28 ligands modelled in this entry, 20 are monoatomic - leaving 8 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	$\# Z > 2$
5	NAG	A	1768	1	14,14,15	0.51	0	15,19,21	1.03	1 (6%)
5	NAG	B	1768	1	14,14,15	0.52	0	15,19,21	1.62	1 (6%)
5	NAG	B	1770	1	14,14,15	0.61	0	15,19,21	2.85	4 (26%)
5	NAG	B	1773	-	14,14,15	0.59	0	15,19,21	1.85	3 (20%)
5	NAG	C	1768	1	14,14,15	0.59	0	15,19,21	0.78	0
5	NAG	D	1768	1	14,14,15	0.56	0	15,19,21	1.36	2 (13%)
5	NAG	D	1770	1	14,14,15	0.65	0	15,19,21	1.33	2 (13%)
5	NAG	D	1774	-	14,14,15	0.61	0	15,19,21	1.73	3 (20%)

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
5	NAG	A	1768	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1768	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1770	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1773	-	-	0/6/23/26	0/1/1/1
5	NAG	C	1768	1	-	0/6/23/26	0/1/1/1
5	NAG	D	1768	1	-	0/6/23/26	0/1/1/1
5	NAG	D	1770	1	-	0/6/23/26	0/1/1/1
5	NAG	D	1774	-	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1770	NAG	C6-C5-C4	-2.74	106.26	113.02

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1773	NAG	C3-C2-N2	2.15	115.72	110.56
5	D	1770	NAG	C3-C4-C5	2.24	114.10	110.20
5	D	1774	NAG	C3-C2-N2	2.28	116.01	110.56
5	D	1768	NAG	C4-C3-C2	2.29	114.79	111.23
5	A	1768	NAG	C1-O5-C5	2.93	115.96	112.25
5	D	1774	NAG	C1-O5-C5	3.39	116.55	112.25
5	B	1773	NAG	C1-O5-C5	3.45	116.63	112.25
5	D	1770	NAG	C4-C3-C2	3.54	116.73	111.23
5	B	1770	NAG	C4-C3-C2	3.66	116.92	111.23
5	B	1770	NAG	C3-C4-C5	3.76	116.75	110.20
5	D	1768	NAG	C1-O5-C5	3.79	117.06	112.25
5	D	1774	NAG	C2-N2-C7	4.28	128.54	123.04
5	B	1773	NAG	C2-N2-C7	5.14	129.64	123.04
5	B	1768	NAG	C1-O5-C5	5.60	119.35	112.25
5	B	1770	NAG	C1-O5-C5	8.75	123.36	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1768	NAG	1	0
5	B	1773	NAG	1	0
5	D	1774	NAG	1	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	705/735 (95%)	-0.34	16 (2%) 64 67	26, 34, 64, 88	0
1	B	708/735 (96%)	-0.35	11 (1%) 74 78	24, 34, 67, 99	0
1	C	705/735 (95%)	-0.35	12 (1%) 73 76	25, 35, 64, 84	0
1	D	708/735 (96%)	-0.32	14 (1%) 68 72	24, 34, 68, 101	0
All	All	2826/2940 (96%)	-0.34	53 (1%) 70 73	24, 34, 67, 101	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	762	HIS	11.7
1	A	203	HIS	7.3
1	C	762	HIS	6.9
1	C	203	HIS	6.5
1	B	204	ARG	5.9
1	D	761	SER	5.9
1	B	762	HIS	5.8
1	A	204	ARG	5.7
1	B	40	HIS	5.5
1	A	747	ALA	4.4
1	C	204	ARG	4.0
1	A	742	LEU	3.9
1	D	735	GLU	3.9
1	C	751	ASP	3.8
1	D	747	ALA	3.8
1	C	761	SER	3.7
1	D	40	HIS	3.6
1	D	204	ARG	3.6
1	D	203	HIS	3.5
1	B	203	HIS	3.2
1	A	741	CYS	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	745	ALA	3.1
1	B	761	SER	3.1
1	A	751	ASP	3.1
1	B	736	VAL	3.1
1	C	745	ALA	3.0
1	C	735	GLU	3.0
1	D	733	ALA	2.9
1	C	747	ALA	2.9
1	B	735	GLU	2.8
1	A	734	CYS	2.7
1	B	503	ALA	2.6
1	A	736	VAL	2.5
1	C	730	ASP	2.5
1	A	205	GLY	2.4
1	C	736	VAL	2.4
1	A	729	GLN	2.4
1	D	734	CYS	2.3
1	D	746	ALA	2.3
1	A	735	GLU	2.3
1	C	728	ASP	2.3
1	D	748	CYS	2.3
1	C	746	ALA	2.2
1	A	738	PRO	2.2
1	D	758	GLY	2.2
1	A	745	ALA	2.1
1	A	746	ALA	2.1
1	D	760	PHE	2.0
1	B	734	CYS	2.0
1	A	737	ASN	2.0
1	B	759	GLY	2.0
1	A	760	PHE	2.0
1	D	504	THR	2.0

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(\AA^2)	Q<0.9
1	TPQ	A	471	14/15	0.86	0.19	-	40,48,51,52	0
1	TPQ	D	471	14/15	0.88	0.18	-	38,46,49,49	0
1	TPQ	B	471	14/15	0.89	0.20	-	39,47,49,50	0
1	TPQ	C	471	14/15	0.91	0.19	-	38,46,49,49	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(\AA^2)	Q<0.9
4	NAG	A	1770	14/15	0.78	0.43	4.69	77,81,81,82	0
4	NAG	A	1767	14/15	0.89	0.22	4.12	58,64,67,73	0
4	NAG	C	1770	14/15	0.83	0.38	3.62	76,79,80,81	0
4	NAG	D	1771	14/15	0.82	0.29	1.68	74,75,77,77	0
4	NAG	C	1767	14/15	0.94	0.16	1.62	58,64,66,72	0
4	NAG	B	1767	14/15	0.91	0.17	0.96	58,63,65,68	0
2	NAG	A	1762	14/15	0.96	0.13	0.57	43,46,49,55	0
4	NAG	D	1767	14/15	0.91	0.17	0.53	59,66,68,71	0
2	NAG	B	1762	14/15	0.96	0.12	0.13	39,43,48,55	0
2	NAG	C	1762	14/15	0.93	0.13	-0.03	43,49,51,55	0
3	NDG	C	1764	14/15	0.95	0.12	-0.36	42,48,53,58	0
9	NDG	B	1764	14/15	0.96	0.12	-0.37	39,45,51,54	0
10	NAG	D	1762	14/15	0.97	0.12	-0.41	40,44,49,55	0
3	NDG	A	1764	14/15	0.95	0.12	-0.45	39,44,48,53	0
9	NDG	D	1764	14/15	0.98	0.10	-1.02	34,43,52,52	0
3	FUL	C	1766	10/11	0.93	0.15	-	47,50,52,53	0
9	MAN	D	1775	11/12	0.73	0.28	-	79,81,82,82	0
2	NAG	C	1763	14/15	0.93	0.12	-	59,61,66,70	0
10	NAG	D	1763	14/15	0.91	0.14	-	61,64,68,74	0
2	NAG	B	1763	14/15	0.87	0.18	-	62,66,70,75	0
2	BMA	C	1769	11/12	0.88	0.23	-	73,75,75,76	0
9	BMA	B	1772	11/12	0.82	0.13	-	69,70,72,75	0
2	NAG	A	1763	14/15	0.89	0.18	-	59,63,65,69	0
3	NAG	C	1765	14/15	0.84	0.25	-	62,65,69,69	0
9	MAN	B	1774	11/12	0.84	0.28	-	78,79,80,81	0
4	NAG	B	1771	14/15	0.82	0.29	-	71,73,73,74	0
2	BMA	B	1769	11/12	0.74	0.18	-	79,82,82,82	0
9	FUL	D	1766	10/11	0.86	0.18	-	52,55,57,58	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	FUL	A	1766	10/11	0.93	0.14	-	50,51,52,52	0
2	BMA	A	1769	11/12	0.84	0.24	-	72,74,75,75	0
9	NAG	D	1765	14/15	0.95	0.14	-	55,59,62,66	0
3	NAG	A	1765	14/15	0.84	0.23	-	57,61,62,62	0
10	BMA	D	1769	11/12	0.85	0.16	-	79,82,86,88	0
9	BMA	D	1772	11/12	0.83	0.14	-	67,70,72,76	0
9	NAG	B	1765	14/15	0.88	0.17	-	54,61,67,68	0
9	FUL	B	1766	10/11	0.90	0.22	-	53,55,56,57	0
10	MAN	D	1773	11/12	0.70	0.20	-	89,90,91,91	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(\AA^2)	Q<0.9
8	CL	C	1774	1/1	1.00	0.10	-0.71	15,15,15,15	0
8	CL	D	1779	1/1	0.99	0.11	-0.94	16,16,16,16	0
8	CL	A	1774	1/1	0.99	0.11	-1.04	19,19,19,19	0
6	CA	B	1777	1/1	0.99	0.07	-2.47	31,31,31,31	0
6	CA	A	1773	1/1	0.97	0.08	-2.65	40,40,40,40	0
6	CA	D	1776	1/1	0.99	0.07	-2.78	25,25,25,25	0
6	CA	C	1773	1/1	0.94	0.07	-2.94	42,42,42,42	0
8	CL	B	1778	1/1	0.99	0.11	-2.98	18,18,18,18	0
6	CA	D	1778	1/1	0.98	0.07	-3.02	29,29,29,29	0
7	CU	C	1772	1/1	0.99	0.06	-3.66	36,36,36,36	0
6	CA	A	1771	1/1	0.99	0.05	-3.95	29,29,29,29	0
6	CA	C	1771	1/1	0.99	0.06	-4.87	28,28,28,28	0
6	CA	B	1775	1/1	0.99	0.04	-6.76	27,27,27,27	0
7	CU	A	1772	1/1	0.99	0.08	-	33,33,33,33	0
5	NAG	A	1768	14/15	0.85	0.23	-	64,69,70,70	0
5	NAG	D	1774	14/15	0.76	0.20	-	83,86,87,87	0
8	CL	C	1775	1/1	0.98	0.09	-	24,24,24,24	1
5	NAG	D	1770	14/15	0.88	0.17	-	60,65,67,67	0
5	NAG	C	1768	14/15	0.88	0.21	-	61,66,66,68	0
8	CL	D	1780	1/1	0.94	0.07	-	21,21,21,21	1
8	CL	A	1775	1/1	0.99	0.06	-	24,24,24,24	1
7	CU	D	1777	1/1	0.99	0.07	-	35,35,35,35	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	NAG	D	1768	14/15	0.87	0.18	-	62,66,69,70	0
5	NAG	B	1773	14/15	0.84	0.17	-	77,78,78,79	0
5	NAG	B	1768	14/15	0.87	0.18	-	62,67,70,71	0
5	NAG	B	1770	14/15	0.84	0.21	-	60,65,66,66	0
7	CU	B	1776	1/1	1.00	0.08	-	35,35,35,35	0
8	CL	B	1779	1/1	0.99	0.07	-	18,18,18,18	1

6.5 Other polymers [i](#)

There are no such residues in this entry.