



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:45 AM GMT

PDB ID : 3PY4
Title : Crystal structure of bovine lactoperoxidase in complex with paracetamol at 2.4Å resolution
Authors : Pandey, N.; Sing, R.P.; Singh, A.K.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2010-12-11
Resolution : 2.42 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

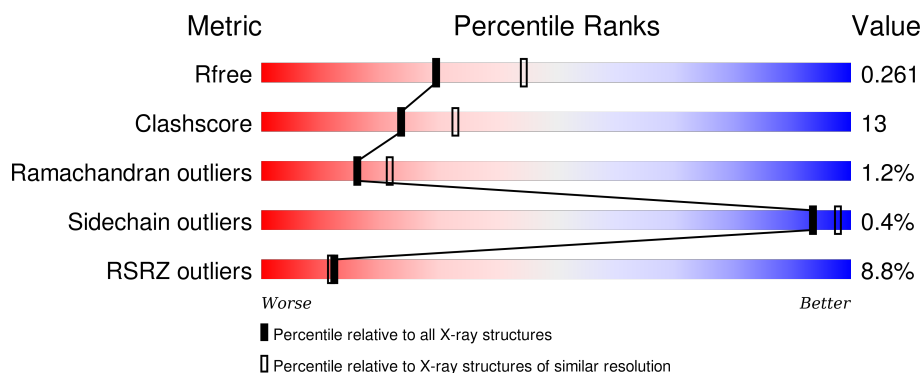
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (2.44-2.40)

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	595	<div> <div>9%</div> <div>79%</div> <div>20%</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
11	EDO	A	702	-	-	X	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	IOD	A	616	-	-	X	X
8	TYL	A	598	-	-	X	-
9	MPD	A	613	-	-	X	X

2 Entry composition [i](#)

There are 13 unique types of molecules in this entry. The entry contains 5150 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 Lactoperoxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	595	Total	C	N	O	P	S	0	0	0
			4774	3037	847	863	1	26			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 6 is a polymer of unknown type called SUGAR (3-MER).

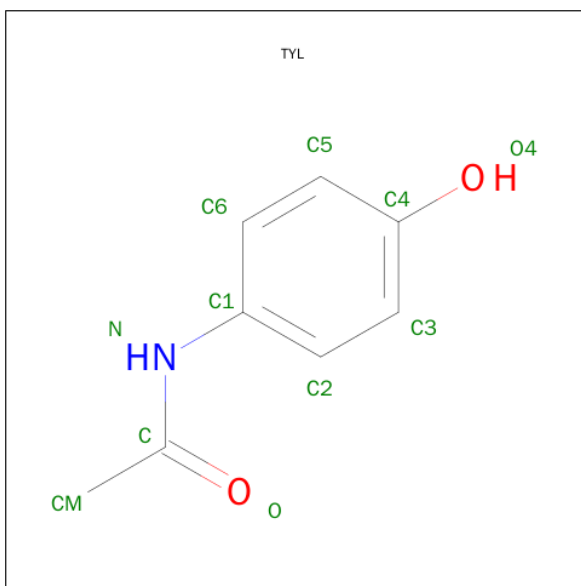
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 7 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	15	Total	I	0	0
			15	15		

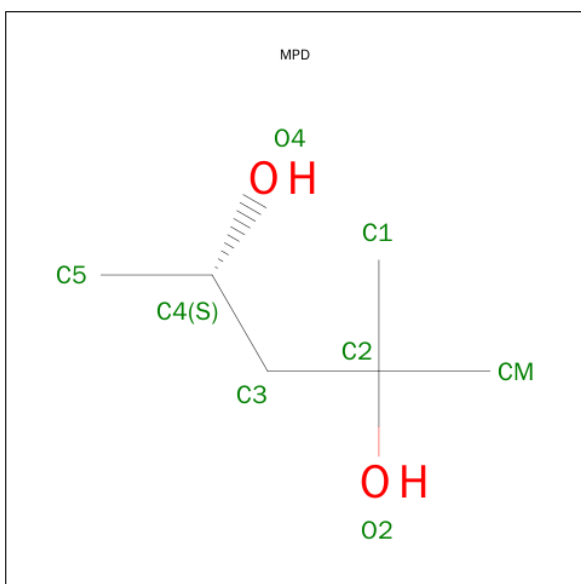
- Molecule 8 is N-(4-HYDROXYPHENYL)ACETAMIDE (TYLENOL) (three-letter code:

TYL) (formula: C₈H₉NO₂).



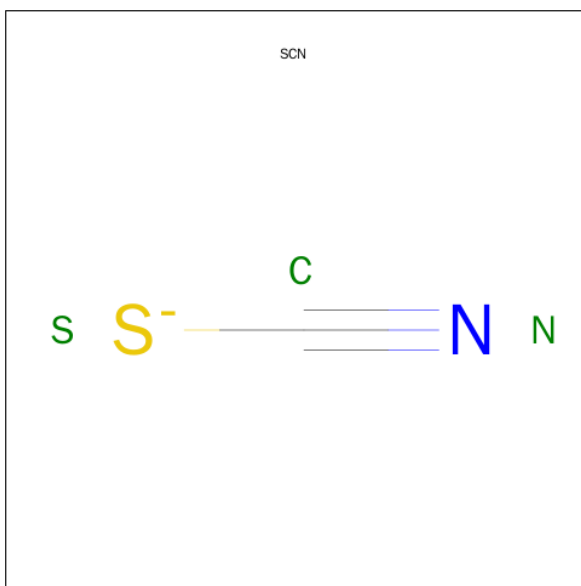
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			11	8	1	2		

- Molecule 9 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



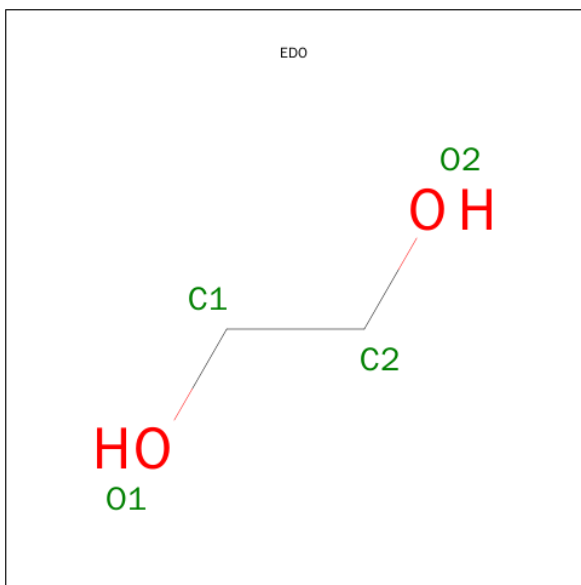
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			8	6	2		

- Molecule 10 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



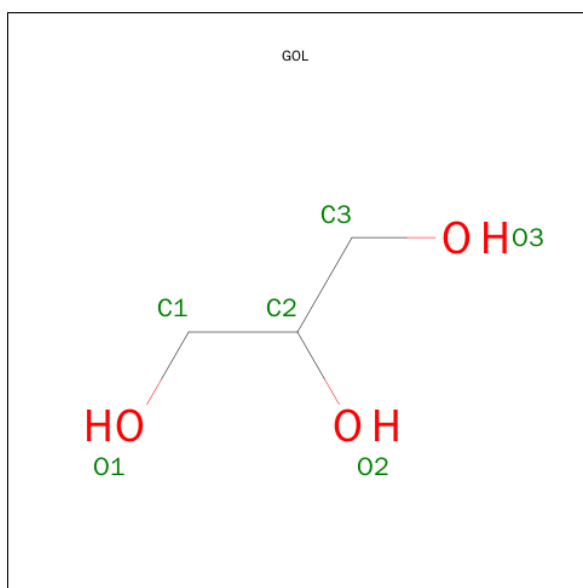
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	A	1	Total	C	N	S	0	0
			3	1	1	1		

- Molecule 11 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 12 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

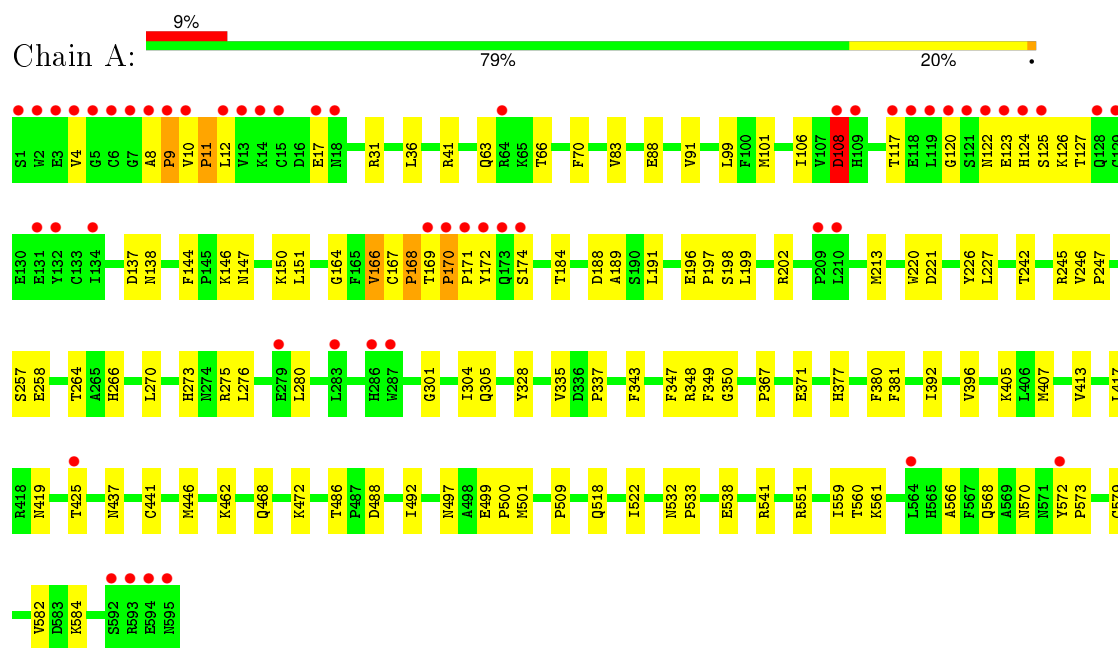
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	186	Total	O	0	0
			186	186		

3 Residue-property plots [i](#)

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: Lactoperoxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.81Å 80.67Å 74.44Å 90.00° 103.07° 90.00°	Depositor
Resolution (Å)	43.90 – 2.42 43.95 – 2.42	Depositor EDS
% Data completeness (in resolution range)	94.0 (43.90-2.42) 93.9 (43.95-2.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.48 (at 2.42Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.225 , 0.258 0.218 , 0.261	Depositor DCC
R_{free} test set	1149 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	45.1	Xtriage
Anisotropy	0.774	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 37.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 22419 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5150	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% 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: TYL, GOL, MPD, SCN, NAG, SEP, CA, EDO, HEM, IOD, 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	0.48	0/4891	0.73	4/6634 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	ASP	CA-CB-CG	12.29	140.45	113.40
1	A	108	ASP	N-CA-CB	9.12	127.01	110.60
1	A	108	ASP	CB-CA-C	-7.17	96.06	110.40
1	A	486	THR	N-CA-CB	5.35	120.47	110.30

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	4774	0	4687	126	0
2	A	1	0	0	0	0
3	A	43	0	30	11	0
4	A	28	0	26	1	0
5	A	28	0	25	1	0
6	A	39	0	34	0	0
7	A	15	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	11	0	8	8	0
9	A	8	0	14	8	0
10	A	3	0	0	0	0
11	A	8	0	12	6	0
12	A	6	0	8	1	0
13	A	186	0	0	7	0
All	All	5150	0	4844	129	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 13.

All (129) 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:258:GLU:OE2	3:A:605:HEM:CMB	1.73	1.35
1:A:108:ASP:OD2	3:A:605:HEM:CMD	1.76	1.34
1:A:258:GLU:OE2	3:A:605:HEM:HMB1	1.06	1.21
1:A:168:PRO:HG3	1:A:172:TYR:HB2	1.24	1.14
1:A:108:ASP:OD2	3:A:605:HEM:HMD1	0.98	1.13
1:A:227:LEU:HD23	1:A:270:LEU:HD22	1.40	1.02
1:A:551:ARG:HD3	1:A:584:LYS:HA	1.40	1.00
1:A:258:GLU:HG3	8:A:598:TYL:H2	1.47	0.96
1:A:170:PRO:HB2	1:A:171:PRO:HD3	1.50	0.93
1:A:197:PRO:HA	11:A:702:EDO:H11	1.55	0.89
1:A:258:GLU:HB2	8:A:598:TYL:H2	1.60	0.84
1:A:197:PRO:HB3	11:A:702:EDO:H22	1.60	0.83
1:A:258:GLU:CG	8:A:598:TYL:H2	2.09	0.81
1:A:108:ASP:OD2	3:A:605:HEM:C2D	2.33	0.81
1:A:197:PRO:CA	11:A:702:EDO:H11	2.11	0.80
1:A:258:GLU:CB	8:A:598:TYL:H2	2.13	0.79
1:A:117:THR:HG22	1:A:164:GLY:HA2	1.63	0.78
1:A:258:GLU:HG3	8:A:598:TYL:C2	2.14	0.77
1:A:168:PRO:HG3	1:A:172:TYR:CB	2.10	0.77
1:A:197:PRO:HA	11:A:702:EDO:C1	2.15	0.77
1:A:196:GLU:HB3	1:A:198:SEP:O2P	1.85	0.76
1:A:335:VAL:O	1:A:337:PRO:HD3	1.88	0.73
1:A:17:GLU:O	1:A:31:ARG:NH2	2.21	0.72
1:A:258:GLU:OE2	3:A:605:HEM:HMB2	1.86	0.72
1:A:197:PRO:HD2	1:A:198:SEP:O3P	1.90	0.71
1:A:9:PRO:O	1:A:11:PRO:HD3	1.92	0.69
1:A:91:VAL:HG22	1:A:405:LYS:HG3	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:ARG:NH1	9:A:613:MPD:H12	2.07	0.68
1:A:499:GLU:OE1	1:A:509:PRO:HD2	1.94	0.66
1:A:407:MET:HB3	1:A:501:MET:CE	2.24	0.66
1:A:99:LEU:HD23	1:A:566:ALA:HB1	1.78	0.65
1:A:227:LEU:HD23	1:A:270:LEU:CD2	2.20	0.64
1:A:328:TYR:OH	1:A:532:ASN:HB2	1.98	0.64
1:A:167:CYS:HB2	1:A:168:PRO:HD3	1.79	0.63
1:A:264:THR:HG23	1:A:392:ILE:HB	1.82	0.61
1:A:197:PRO:HB3	11:A:702:EDO:C2	2.31	0.60
1:A:348:ARG:HH11	1:A:437:ASN:ND2	1.98	0.60
1:A:349:PHE:CB	1:A:497:ASN:HD21	2.15	0.60
1:A:150:LYS:HE2	1:A:419:ASN:OD1	2.01	0.59
1:A:276:LEU:O	1:A:280:LEU:HG	2.03	0.59
1:A:213:MET:HG2	1:A:273:HIS:CD2	2.39	0.58
1:A:174:SER:HA	13:A:812:HOH:O	2.04	0.58
1:A:301:GLY:O	1:A:305:GLN:HG3	2.04	0.57
1:A:348:ARG:HH11	1:A:437:ASN:HD22	1.51	0.57
8:A:598:TYL:H5	13:A:797:HOH:O	2.05	0.57
1:A:441:CYS:SG	1:A:446:MET:HG3	2.46	0.56
1:A:425:THR:O	1:A:425:THR:HG22	2.07	0.55
1:A:381:PHE:CE1	8:A:598:TYL:HM3	2.42	0.55
1:A:275:ARG:HD2	13:A:723:HOH:O	2.08	0.53
1:A:258:GLU:OE2	3:A:605:HEM:C2B	2.57	0.53
1:A:91:VAL:HG21	1:A:405:LYS:HE2	1.91	0.53
1:A:202:ARG:NH1	9:A:613:MPD:C1	2.71	0.53
1:A:170:PRO:HB2	1:A:171:PRO:CD	2.31	0.52
1:A:198:SEP:H	1:A:198:SEP:P	2.32	0.52
4:A:604:NAG:H61	13:A:815:HOH:O	2.10	0.52
1:A:8:ALA:N	1:A:9:PRO:HD2	2.25	0.52
1:A:579:CYS:O	1:A:582:VAL:HB	2.10	0.52
1:A:441:CYS:SG	1:A:492:ILE:HG22	2.50	0.51
1:A:227:LEU:HD11	1:A:266:HIS:HB3	1.92	0.51
1:A:184:THR:OG1	1:A:188:ASP:OD2	2.29	0.50
1:A:63:GLN:H	1:A:63:GLN:CD	2.14	0.50
1:A:348:ARG:NH2	3:A:605:HEM:HAD1	2.27	0.50
1:A:199:LEU:HB2	9:A:613:MPD:HM1	1.92	0.50
1:A:551:ARG:HD2	1:A:582:VAL:HG12	1.94	0.49
1:A:189:ALA:HB2	1:A:304:ILE:HD12	1.95	0.49
1:A:371:GLU:OE2	13:A:804:HOH:O	2.20	0.49
1:A:568:GLN:HE21	1:A:570:ASN:ND2	2.11	0.48
1:A:392:ILE:O	1:A:396:VAL:HG23	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:VAL:HG12	1:A:413:VAL:HB	1.95	0.48
1:A:561:LYS:HD3	1:A:572:TYR:CE1	2.48	0.47
1:A:257:SER:O	1:A:381:PHE:HA	2.14	0.47
1:A:532:ASN:OD1	1:A:533:PRO:HD2	2.14	0.47
1:A:220:TRP:HB2	13:A:690:HOH:O	2.15	0.47
1:A:106:ILE:HG23	1:A:191:LEU:HD11	1.97	0.47
1:A:91:VAL:HG11	7:A:616:IOD:I	2.85	0.47
1:A:538:GLU:OE2	1:A:541:ARG:NH2	2.48	0.47
1:A:4:VAL:O	1:A:4:VAL:HG13	2.15	0.47
1:A:11:PRO:HD2	1:A:12:LEU:H	1.80	0.46
1:A:166:VAL:HG23	1:A:167:CYS:H	1.80	0.46
1:A:560:THR:HA	1:A:579:CYS:SG	2.56	0.46
1:A:472:LYS:HD2	1:A:500:PRO:HG2	1.97	0.46
1:A:561:LYS:HD3	1:A:572:TYR:HE1	1.81	0.46
1:A:246:VAL:HA	1:A:247:PRO:HD3	1.79	0.46
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.51	0.46
1:A:199:LEU:HD13	9:A:613:MPD:HM3	1.98	0.46
1:A:88:GLU:OE1	7:A:616:IOD:I	3.04	0.45
1:A:99:LEU:CD2	1:A:566:ALA:HB1	2.46	0.45
1:A:572:TYR:CG	1:A:573:PRO:HA	2.51	0.45
1:A:202:ARG:HH12	9:A:613:MPD:H12	1.81	0.45
1:A:462:LYS:NZ	1:A:488:ASP:OD1	2.46	0.45
1:A:349:PHE:HB2	1:A:497:ASN:HD21	1.81	0.45
1:A:196:GLU:HA	1:A:197:PRO:HD3	1.82	0.45
1:A:122:ASN:ND2	13:A:772:HOH:O	2.49	0.45
1:A:221:ASP:HB2	1:A:226:TYR:CE2	2.51	0.45
1:A:350:GLY:HA3	3:A:605:HEM:CBC	2.47	0.44
1:A:343:PHE:CG	1:A:518:GLN:HG2	2.52	0.44
1:A:10:VAL:HB	1:A:41:ARG:HE	1.83	0.44
1:A:91:VAL:CG1	7:A:616:IOD:I	3.37	0.43
1:A:120:GLY:O	1:A:126:LYS:HD3	2.18	0.43
1:A:381:PHE:CZ	8:A:598:TYL:HM3	2.54	0.43
1:A:213:MET:HB3	1:A:270:LEU:HD11	2.01	0.43
5:A:599:NAG:H61	5:A:600:NAG:O7	2.19	0.42
1:A:167:CYS:HB2	1:A:168:PRO:CD	2.46	0.42
1:A:213:MET:CB	1:A:270:LEU:HD11	2.49	0.42
1:A:468:GLN:O	1:A:472:LYS:N	2.52	0.42
1:A:101:MET:SD	1:A:101:MET:C	2.98	0.42
1:A:349:PHE:CD1	1:A:349:PHE:C	2.93	0.42
1:A:202:ARG:HH11	9:A:613:MPD:C1	2.33	0.42
1:A:242:THR:O	1:A:245:ARG:NH1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:ARG:HD3	9:A:613:MPD:O2	2.20	0.42
1:A:522:ILE:HD12	1:A:522:ILE:C	2.40	0.42
1:A:559:ILE:HG23	7:A:612:IOD:I	2.90	0.41
1:A:123:GLU:HG3	1:A:125:SER:H	1.85	0.41
1:A:197:PRO:CB	11:A:702:EDO:H11	2.50	0.41
1:A:572:TYR:CD1	1:A:573:PRO:HA	2.56	0.41
1:A:146:LYS:NZ	1:A:147:ASN:HD21	2.18	0.41
1:A:188:ASP:O	1:A:189:ALA:HB3	2.20	0.41
1:A:11:PRO:CD	1:A:12:LEU:H	2.33	0.41
1:A:66:THR:HB	1:A:70:PHE:N	2.35	0.41
1:A:417:LEU:CD2	3:A:605:HEM:HMB3	2.51	0.41
1:A:170:PRO:CB	1:A:171:PRO:HD3	2.34	0.41
1:A:137:ASP:HB3	1:A:138:ASN:H	1.65	0.41
1:A:568:GLN:HE21	1:A:570:ASN:HD21	1.69	0.40
1:A:221:ASP:HB2	1:A:226:TYR:CZ	2.56	0.40
1:A:144:PHE:HB2	1:A:151:LEU:HD13	2.02	0.40
1:A:124:HIS:O	1:A:127:THR:N	2.54	0.40
1:A:417:LEU:HD21	3:A:605:HEM:HMB3	2.04	0.40
1:A:202:ARG:HH11	9:A:613:MPD:H12	1.81	0.40
1:A:36:LEU:HD11	12:A:617:GOL:H11	2.02	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	592/595 (100%)	543 (92%)	42 (7%)	7 (1%)	16 22

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	VAL

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Mol	Chain	Res	Type
1	A	169	THR
1	A	170	PRO
1	A	11	PRO
1	A	9	PRO
1	A	168	PRO
1	A	367	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	517/517 (100%)	515 (100%)	2 (0%)	93	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	ASP
1	A	347	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	A	147	ASN
1	A	437	ASN
1	A	468	GLN
1	A	497	ASN
1	A	570	ASN

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	SEP	A	198	1	8,9,10	1.53	1 (12%)	8,12,14	1.54	2 (25%)

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	SEP	A	198	1	-	0/6/8/10	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	198	SEP	P-O1P	3.14	1.61	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	198	SEP	O-C-CA	-2.13	119.94	125.49
1	A	198	SEP	OG-CB-CA	3.31	111.09	108.27

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	198	SEP	3	0

5.5 Carbohydrates ⓘ

5 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	$\# Z > 2$
5	NAG	A	599	1,5	14,14,15	0.66	0	15,19,21	1.21	3 (20%)
5	NAG	A	600	5	14,14,15	0.66	0	15,19,21	1.24	2 (13%)
6	NAG	A	601	1,6	14,14,15	0.95	1 (7%)	15,19,21	1.08	2 (13%)
6	NAG	A	602	6	14,14,15	0.60	0	15,19,21	1.73	2 (13%)
6	MAN	A	603	6	11,11,12	0.63	0	14,15,17	1.33	2 (14%)

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	599	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	600	5	-	0/6/23/26	0/1/1/1
6	NAG	A	601	1,6	-	0/6/23/26	0/1/1/1
6	NAG	A	602	6	-	0/6/23/26	0/1/1/1
6	MAN	A	603	6	-	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)
6	A	601	NAG	O5-C1	-2.27	1.39	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	601	NAG	C1-O5-C5	-2.51	109.07	112.25
5	A	599	NAG	C3-C2-N2	-2.35	104.94	110.56
6	A	601	NAG	C2-N2-C7	-2.15	120.27	123.04
5	A	599	NAG	C1-O5-C5	-2.15	109.52	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	599	NAG	O4-C4-C3	-2.09	105.63	110.34
6	A	602	NAG	C4-C3-C2	-2.01	108.11	111.23
5	A	600	NAG	C2-N2-C7	2.17	125.83	123.04
5	A	600	NAG	C4-C3-C2	2.64	115.33	111.23
6	A	603	MAN	C1-O5-C5	3.04	116.10	112.25
6	A	603	MAN	C1-C2-C3	3.21	113.34	109.54
6	A	602	NAG	C1-O5-C5	4.86	118.42	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	599	NAG	1	0
5	A	600	NAG	1	0

5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 16 are monoatomic - leaving 9 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	$\# Z > 2$
4	NAG	A	596	1	14,14,15	0.83	1 (7%)	15,19,21	1.66	4 (26%)
8	TYL	A	598	-	11,11,11	1.11	1 (9%)	14,14,14	1.42	1 (7%)
4	NAG	A	604	1	14,14,15	0.54	0	15,19,21	1.40	1 (6%)
3	HEM	A	605	1	30,50,50	2.88	13 (43%)	24,82,82	2.27	8 (33%)
9	MPD	A	613	-	6,7,7	0.25	0	7,10,10	0.47	0
12	GOL	A	617	-	5,5,5	0.47	0	5,5,5	0.14	0
10	SCN	A	700	-	2,2,2	2.40	1 (50%)	1,1,1	0.07	0
11	EDO	A	701	-	3,3,3	0.70	0	2,2,2	0.40	0
11	EDO	A	702	-	3,3,3	1.04	0	2,2,2	0.76	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
4	NAG	A	596	1	-	0/6/23/26	0/1/1/1
8	TYL	A	598	-	-	2/4/4/4	0/1/1/1
4	NAG	A	604	1	-	0/6/23/26	0/1/1/1
3	HEM	A	605	1	-	0/10/54/54	0/0/8/8
9	MPD	A	613	-	-	0/5/5/5	0/0/0/0
12	GOL	A	617	-	-	0/4/4/4	0/0/0/0
10	SCN	A	700	-	-	0/0/0/0	0/0/0/0
11	EDO	A	701	-	-	0/1/1/1	0/0/0/0
11	EDO	A	702	-	-	0/1/1/1	0/0/0/0

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	605	HEM	C3B-C4B	-11.26	1.41	1.51
8	A	598	TYL	C1-N	-3.28	1.35	1.41
3	A	605	HEM	C2D-C1D	-3.16	1.41	1.51
3	A	605	HEM	C3D-C4D	-3.05	1.47	1.51
3	A	605	HEM	C1A-CHA	-2.98	1.31	1.39
3	A	605	HEM	C2B-C1B	-2.82	1.42	1.51
3	A	605	HEM	C2C-C1C	-2.71	1.47	1.52
3	A	605	HEM	C2D-C3D	-2.54	1.47	1.54
4	A	596	NAG	O5-C1	-2.10	1.40	1.43
3	A	605	HEM	CMC-C2C	2.45	1.58	1.53
3	A	605	HEM	FE-NC	2.49	2.05	1.95
3	A	605	HEM	C3B-CAB	2.88	1.56	1.51
3	A	605	HEM	CAD-C3D	3.06	1.60	1.54
10	A	700	SCN	C-S	3.28	1.84	1.63
3	A	605	HEM	CAA-C2A	3.34	1.57	1.52
3	A	605	HEM	C3C-CAC	3.72	1.58	1.51

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	598	TYL	C1-N-C	-5.18	120.00	128.03
3	A	605	HEM	CAA-C2A-C1A	-2.43	124.37	127.01
3	A	605	HEM	C1D-CHD-C4C	-2.13	122.26	125.82
4	A	596	NAG	C2-N2-C7	2.10	125.74	123.04
3	A	605	HEM	C3C-CAC-CBC	2.31	127.99	124.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	605	HEM	CAD-C3D-C2D	2.92	121.60	113.22
4	A	596	NAG	C3-C4-C5	2.92	115.28	110.20
3	A	605	HEM	C4B-CHC-C1C	3.03	130.89	125.82
4	A	596	NAG	C4-C3-C2	3.05	115.97	111.23
4	A	596	NAG	C1-O5-C5	3.11	116.19	112.25
3	A	605	HEM	CMC-C2C-C3C	3.69	125.75	116.53
3	A	605	HEM	CMB-C2B-C3B	4.45	127.64	116.53
4	A	604	NAG	C1-O5-C5	4.91	118.48	112.25
3	A	605	HEM	CAD-C3D-C4D	6.56	135.62	112.47

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	598	TYL	O-C-N-C1
8	A	598	TYL	CM-C-N-C1

There are no ring outliers.

6 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	598	TYL	8	0
4	A	604	NAG	1	0
3	A	605	HEM	11	0
9	A	613	MPD	8	0
12	A	617	GOL	1	0
11	A	702	EDO	6	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	594/595 (99%)	0.40	52 (8%) 12 12	31, 51, 93, 120	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	TRP	15.3
1	A	173	GLN	13.0
1	A	121	SER	11.8
1	A	1	SER	10.0
1	A	4	VAL	8.6
1	A	595	ASN	8.4
1	A	13	VAL	6.8
1	A	172	TYR	6.8
1	A	122	ASN	6.7
1	A	120	GLY	6.5
1	A	7	GLY	6.3
1	A	124	HIS	5.8
1	A	174	SER	5.7
1	A	594	GLU	5.6
1	A	12	LEU	5.5
1	A	132	TYR	5.3
1	A	10	VAL	5.3
1	A	5	GLY	5.2
1	A	6	CYS	4.8
1	A	3	GLU	4.3
1	A	593	ARG	4.1
1	A	128	GLN	4.0
1	A	8	ALA	4.0
1	A	592	SER	4.0
1	A	169	THR	3.9
1	A	209	PRO	3.9
1	A	170	PRO	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	125	SER	3.5
1	A	171	PRO	3.3
1	A	17	GLU	3.1
1	A	119	LEU	3.0
1	A	118	GLU	3.0
1	A	64	ARG	3.0
1	A	564	LEU	2.9
1	A	18	ASN	2.8
1	A	283	LEU	2.8
1	A	425	THR	2.7
1	A	134	ILE	2.7
1	A	287	TRP	2.7
1	A	286	HIS	2.6
1	A	9	PRO	2.5
1	A	117	THR	2.4
1	A	14	LYS	2.3
1	A	129	CYS	2.3
1	A	210	LEU	2.3
1	A	108	ASP	2.2
1	A	123	GLU	2.1
1	A	279	GLU	2.1
1	A	109	HIS	2.1
1	A	15	CYS	2.0
1	A	131	GLU	2.0
1	A	572	TYR	2.0

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

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	SEP	A	198	10/11	0.79	0.21	-	49,51,55,57	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
6	NAG	A	601	14/15	0.91	0.14	-0.16	61,64,68,72	0
5	NAG	A	599	14/15	0.93	0.11	-1.39	66,68,72,80	0
5	NAG	A	600	14/15	0.83	0.23	-	89,92,96,99	0
6	MAN	A	603	11/12	0.83	0.33	-	88,90,91,92	0
6	NAG	A	602	14/15	0.91	0.21	-	75,78,82,85	0

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(Å ²)	Q<0.9
9	MPD	A	613	8/8	0.81	0.72	22.97	64,69,71,71	8
11	EDO	A	702	4/4	0.93	0.26	3.58	27,30,31,31	0
7	IOD	A	616	1/1	0.58	0.26	2.69	100,100,100,100	1
8	TYL	A	598	11/11	0.83	0.30	1.60	22,35,37,40	0
10	SCN	A	700	3/3	0.94	0.20	0.56	53,53,54,56	0
3	HEM	A	605	43/43	0.97	0.19	-0.26	28,34,39,44	0
7	IOD	A	623	1/1	0.94	0.10	-0.92	66,66,66,66	1
2	CA	A	606	1/1	0.98	0.14	-1.29	42,42,42,42	0
7	IOD	A	612	1/1	0.99	0.07	-1.50	60,60,60,60	0
7	IOD	A	622	1/1	0.99	0.03	-2.16	74,74,74,74	0
7	IOD	A	621	1/1	0.96	0.07	-2.18	73,73,73,73	1
7	IOD	A	611	1/1	0.99	0.07	-2.45	57,57,57,57	0
7	IOD	A	610	1/1	0.98	0.05	-3.12	71,71,71,71	0
7	IOD	A	607	1/1	0.99	0.08	-3.64	54,54,54,54	0
12	GOL	A	617	6/6	0.92	0.13	-	66,66,67,67	0
4	NAG	A	596	14/15	0.76	0.41	-	68,74,75,77	0
7	IOD	A	625	1/1	0.90	0.11	-	91,91,91,91	1
7	IOD	A	597	1/1	0.98	0.07	-	79,79,79,79	1
7	IOD	A	620	1/1	0.98	0.05	-	80,80,80,80	1
11	EDO	A	701	4/4	0.79	0.29	-	68,68,69,70	0
7	IOD	A	608	1/1	0.99	0.04	-	69,69,69,69	0
4	NAG	A	604	14/15	0.90	0.18	-	71,74,76,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	IOD	A	614	1/1	0.95	0.07	-	69,69,69,69	1
7	IOD	A	624	1/1	0.96	0.05	-	100,100,100,100	0
7	IOD	A	609	1/1	0.84	0.11	-	86,86,86,86	1

6.5 Other polymers [i](#)

There are no such residues in this entry.