



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 29, 2026 – 10:25 am BST

PDB ID : 29TO / pdb_000029to
Title : Structure of monomerized and de-glycosylated native leukocyte myeloperoxidase in complex with the Staphylococcal Peroxidase Inhibitor SPIN
Authors : Leitgeb, U.; Pfanzagl, V.
Deposited on : 2026-04-08
Resolution : 1.44 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

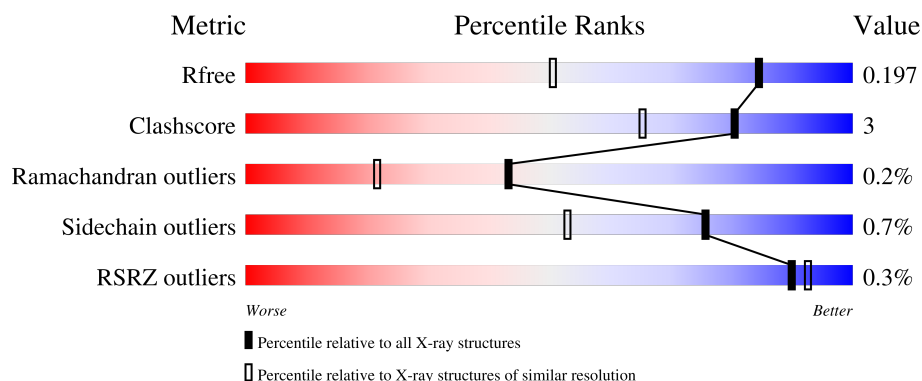
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.44 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3234 (1.46-1.42)
Clashscore	190562	3289 (1.46-1.42)
Ramachandran outliers	187476	3248 (1.46-1.42)
Sidechain outliers	187428	3248 (1.46-1.42)
RSRZ outliers	180081	3234 (1.46-1.42)

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

Mol	Chain	Length	Quality of chain
1	A	114	2% 79% 8% 12%
2	B	467	91% 7%
3	C	73	86% 8% 5%

2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 10793 atoms, of which 5058 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 Myeloperoxidase light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	A	100	Total	C	H	N	O	S	0	1	0
			1580	515	769	146	146	4			

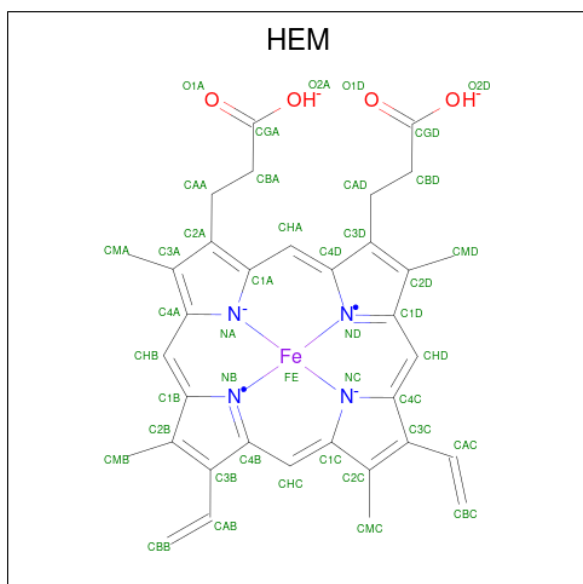
- Molecule 2 is a protein called Myeloperoxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	460	Total	C	H	N	O	S	0	6	0
			7429	2355	3694	690	661	29			

- Molecule 3 is a protein called Myeloperoxidase inhibitor SPIN.

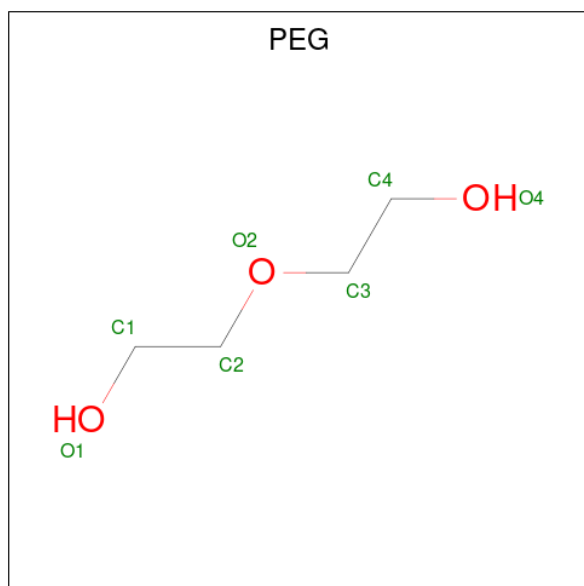
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	69	Total	C	H	N	O	0	2	0
			1113	359	547	95	112			

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	Fe	H	N	O	0	0
			59	34	1	16	4	4		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).

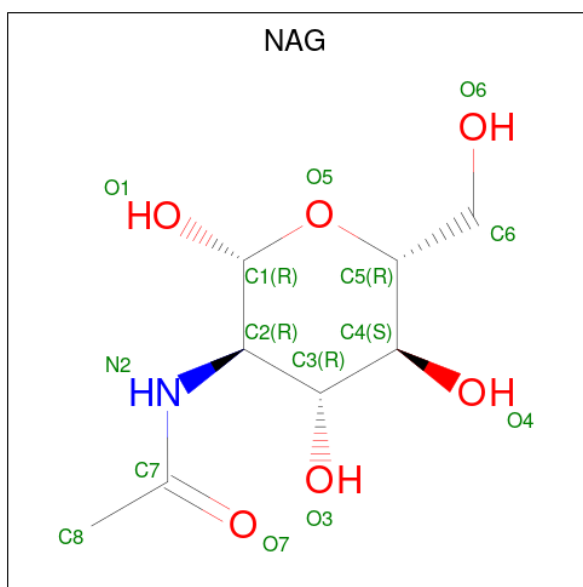


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			15	4	8	3		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		
6	B	8	Total	Cl	0	0
			8	8		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	H	N	O	0	0
			22	8	8	1	5		
7	B	1	Total	C	H	N	O	0	0
			22	8	8	1	5		
7	B	1	Total	C	H	N	O	0	0
			22	8	8	1	5		

- Molecule 8 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Ca	0	0
			1	1		

- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	1	Total	Mg	0	0
			1	1		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	85	Total	O	0	0
			85	85		
10	B	378	Total	O	0	0
			378	378		

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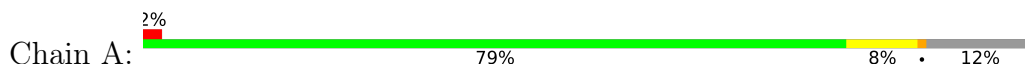
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	C	57	Total	O	0	0
			57	57		

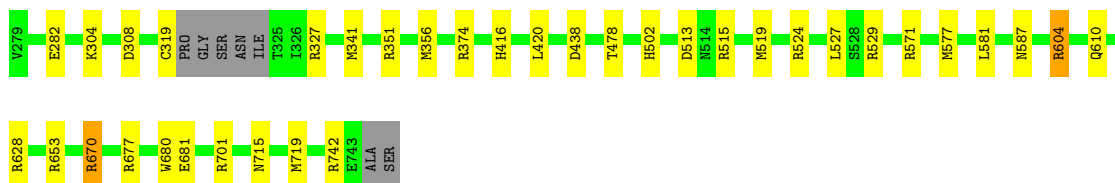
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

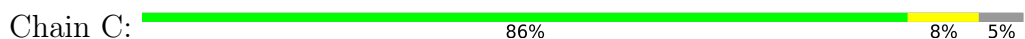
- Molecule 1: Myeloperoxidase light chain



- Molecule 2: Myeloperoxidase



- Molecule 3: Myeloperoxidase inhibitor SPIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.82Å 84.08Å 128.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.38 – 1.44 70.38 – 1.44	Depositor EDS
% Data completeness (in resolution range)	78.2 (70.38-1.44) 78.2 (70.38-1.44)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 1.44Å)	Xtriage
Refinement program	REFMAC 5.8.0431 (refmacat 0.4.126)	Depositor
R, R_{free}	0.140 , 0.196 0.140 , 0.197	Depositor DCC
R_{free} test set	4408 reflections (3.91%)	wwPDB-VP
Wilson B-factor (Å ²)	10.8	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10793	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CSO, HEM, CL, CA, NAG, MG, PEG, YCM

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	0.78	0/838	1.21	5/1140 (0.4%)
2	B	0.77	1/3819 (0.0%)	1.12	10/5173 (0.2%)
3	C	0.76	0/581	1.11	1/780 (0.1%)
All	All	0.77	1/5238 (0.0%)	1.14	16/7093 (0.2%)

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	A	0	3
2	B	0	10
All	All	0	13

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	670	ARG	NE-CZ	5.16	1.38	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	604	ARG	CD-NE-CZ	8.20	135.87	124.40
1	A	239	THR	CA-CB-OG1	8.18	121.86	109.60
1	A	262	ASP	CA-CB-CG	7.79	120.39	112.60
2	B	604	ARG	NE-CZ-NH2	-6.90	112.99	119.20
2	B	604	ARG	NE-CZ-NH1	6.24	127.74	121.50
2	B	478	THR	CA-CB-OG1	-6.20	100.31	109.60
3	C	75	ASP	CB-CA-C	-5.90	97.88	110.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	327	ARG	NE-CZ-NH2	-5.78	114.00	119.20
2	B	438	ASP	CA-CB-CG	5.72	118.32	112.60
1	A	240	ASP	CA-CB-CG	5.63	118.23	112.60
2	B	715	ASN	CA-C-N	5.59	131.77	121.70
2	B	715	ASN	C-N-CA	5.59	131.77	121.70
1	A	264	ASP	CA-CB-CG	5.24	117.84	112.60
2	B	308	ASP	CA-CB-CG	5.22	117.82	112.60
2	B	715	ASN	N-CA-C	-5.20	102.35	110.10
1	A	204	GLU	N-CA-CB	-5.19	102.49	110.12

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	183	ARG	Sidechain
1	A	219[A]	ARG	Sidechain
1	A	219[B]	ARG	Sidechain
2	B	351	ARG	Sidechain
2	B	524	ARG	Sidechain
2	B	571[A]	ARG	Sidechain
2	B	571[B]	ARG	Sidechain
2	B	604	ARG	Sidechain
2	B	628	ARG	Sidechain
2	B	653	ARG	Sidechain
2	B	701	ARG	Sidechain
2	B	742[A]	ARG	Sidechain
2	B	742[B]	ARG	Sidechain

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	811	769	785	2	0
2	B	3735	3694	3752	22	0
3	C	566	547	558	3	0
4	A	43	16	30	0	0
5	A	7	8	10	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	1	0	0	0	0
6	B	8	0	0	0	0
7	B	42	24	39	0	0
8	B	1	0	0	0	0
9	B	1	0	0	0	0
10	A	85	0	0	2	0
10	B	378	0	0	11	0
10	C	57	0	0	0	0
All	All	5735	5058	5174	27	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 3.

All (27) 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)
2:B:319:YCM:HD3	10:B:1170:HOH:O	1.80	0.80
2:B:502:HIS:HD1	2:B:587:ASN:HD21	1.40	0.70
2:B:610:GLN:HG3	10:B:902:HOH:O	1.94	0.66
2:B:282:GLU:OE2	2:B:577[A]:MET:HE2	1.98	0.63
1:A:177:THR:OG1	1:A:179:MET:HE3	2.00	0.62
2:B:319:YCM:NZ2	10:B:903:HOH:O	2.32	0.58
10:A:422:HOH:O	2:B:416:HIS:HE1	1.88	0.57
2:B:356:MET:HA	2:B:356:MET:HE2	1.87	0.55
2:B:529:ARG:NH2	10:B:910:HOH:O	2.39	0.55
5:A:302:PEG:H41	2:B:677:ARG:HG3	1.89	0.54
2:B:680:TRP:CE2	2:B:681:GLU:HG3	2.44	0.53
2:B:519[A]:MET:HE2	10:B:1089:HOH:O	2.09	0.52
10:A:483:HOH:O	2:B:719:MET:HE1	2.10	0.51
2:B:670:ARG:HD3	10:B:1154:HOH:O	2.09	0.51
1:A:183:ARG:HE	5:A:302:PEG:H12	1.76	0.51
2:B:519[A]:MET:CE	10:B:1089:HOH:O	2.61	0.49
10:B:1228:HOH:O	3:C:68:ASN:HB3	2.14	0.47
2:B:513:ASP:HA	2:B:519[B]:MET:HE3	1.97	0.47
2:B:577[A]:MET:HE1	2:B:581:LEU:HD21	1.97	0.47
3:C:57:LEU:HD21	3:C:81:ILE:HG21	1.97	0.46
2:B:304:LYS:HA	2:B:304:LYS:HD3	1.84	0.44
2:B:319:YCM:C	10:B:1234:HOH:O	2.67	0.43
2:B:319:YCM:C	10:B:975:HOH:O	2.68	0.42
3:C:57:LEU:HD12	3:C:98:TYR:CD1	2.55	0.41
2:B:513:ASP:OD2	2:B:515:ARG:NH2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:374[B]:ARG:HD2	10:B:921:HOH:O	2.21	0.41
5:A:302:PEG:H31	2:B:677:ARG:HD2	2.03	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	99/114 (87%)	96 (97%)	2 (2%)	1 (1%)	12	2
2	B	461/467 (99%)	454 (98%)	7 (2%)	0	100	100
3	C	69/73 (94%)	69 (100%)	0	0	100	100
All	All	629/654 (96%)	619 (98%)	9 (1%)	1 (0%)	43	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	208	SER

5.3.2 Protein sidechains [i](#)

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	86/97 (89%)	86 (100%)	0	100	100
2	B	411/410 (100%)	408 (99%)	3 (1%)	76	53

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	63/65 (97%)	62 (98%)	1 (2%)	55	21
All	All	560/572 (98%)	556 (99%)	4 (1%)	76	53

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	341	MET
2	B	420	LEU
2	B	527	LEU
3	C	59	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	220	ASN
2	B	287	GLN
2	B	299	ASN
2	B	416	HIS
2	B	492	ASN
2	B	639	ASN
2	B	706	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	YCM	B	319	2	7,9,10	0.51	0	4,10,12	0.67	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	CSO	B	316	2	3,6,7	1.01	0	0,6,8	-	-

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	YCM	B	319	2	-	1/6/8/10	-
2	CSO	B	316	2	-	0/1/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	319	YCM	CE-CD-SG-CB

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	319	YCM	4	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 11 are monoatomic - leaving 5 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
7	NAG	B	802	2	14,14,15	0.48	0	17,19,21	1.09	2 (11%)
7	NAG	B	801	2	14,14,15	0.64	0	17,19,21	0.94	1 (5%)
7	NAG	B	803	2	14,14,15	0.45	0	17,19,21	1.45	2 (11%)
4	HEM	A	301	1,10,2	50,50,50	1.83	7 (14%)	66,82,82	1.12	7 (10%)
5	PEG	A	302	-	6,6,6	0.30	0	5,5,5	0.29	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
7	NAG	B	802	2	-	1/6/23/26	0/1/1/1
7	NAG	B	801	2	-	0/6/23/26	0/1/1/1
7	NAG	B	803	2	-	2/6/23/26	0/1/1/1
4	HEM	A	301	1,10,2	-	4/14/54/54	-
5	PEG	A	302	-	-	2/4/4/4	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	HEM	FE-NB	5.52	2.12	1.94
4	A	301	HEM	FE-NA	4.88	2.11	1.95
4	A	301	HEM	FE-ND	4.58	2.09	1.94
4	A	301	HEM	FE-NC	3.89	2.08	1.95
4	A	301	HEM	C3C-C4C	-3.05	1.40	1.46
4	A	301	HEM	C3B-C4B	-2.22	1.40	1.44
4	A	301	HEM	C3C-C2C	2.17	1.41	1.37

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	803	NAG	C1-O5-C5	4.31	118.03	112.19
7	B	803	NAG	C1-C2-N2	3.22	115.99	110.49
7	B	802	NAG	C1-C2-N2	-2.75	105.78	110.49
4	A	301	HEM	CHD-C1D-ND	-2.74	121.44	124.42
4	A	301	HEM	CHC-C1C-NC	-2.65	121.57	124.44
4	A	301	HEM	C1B-NB-C4B	2.54	107.70	105.07
4	A	301	HEM	CHD-C1D-C2D	2.45	128.80	124.98
4	A	301	HEM	O1D-CGD-CBD	-2.38	115.43	123.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	301	HEM	O2D-CGD-CBD	2.36	121.61	114.03
7	B	801	NAG	C4-C3-C2	-2.16	107.85	111.02
7	B	802	NAG	O5-C1-C2	-2.13	107.93	111.29
4	A	301	HEM	C4C-CHD-C1D	2.09	130.57	126.06

There are no chirality outliers.

All (9) torsion outliers are listed below:

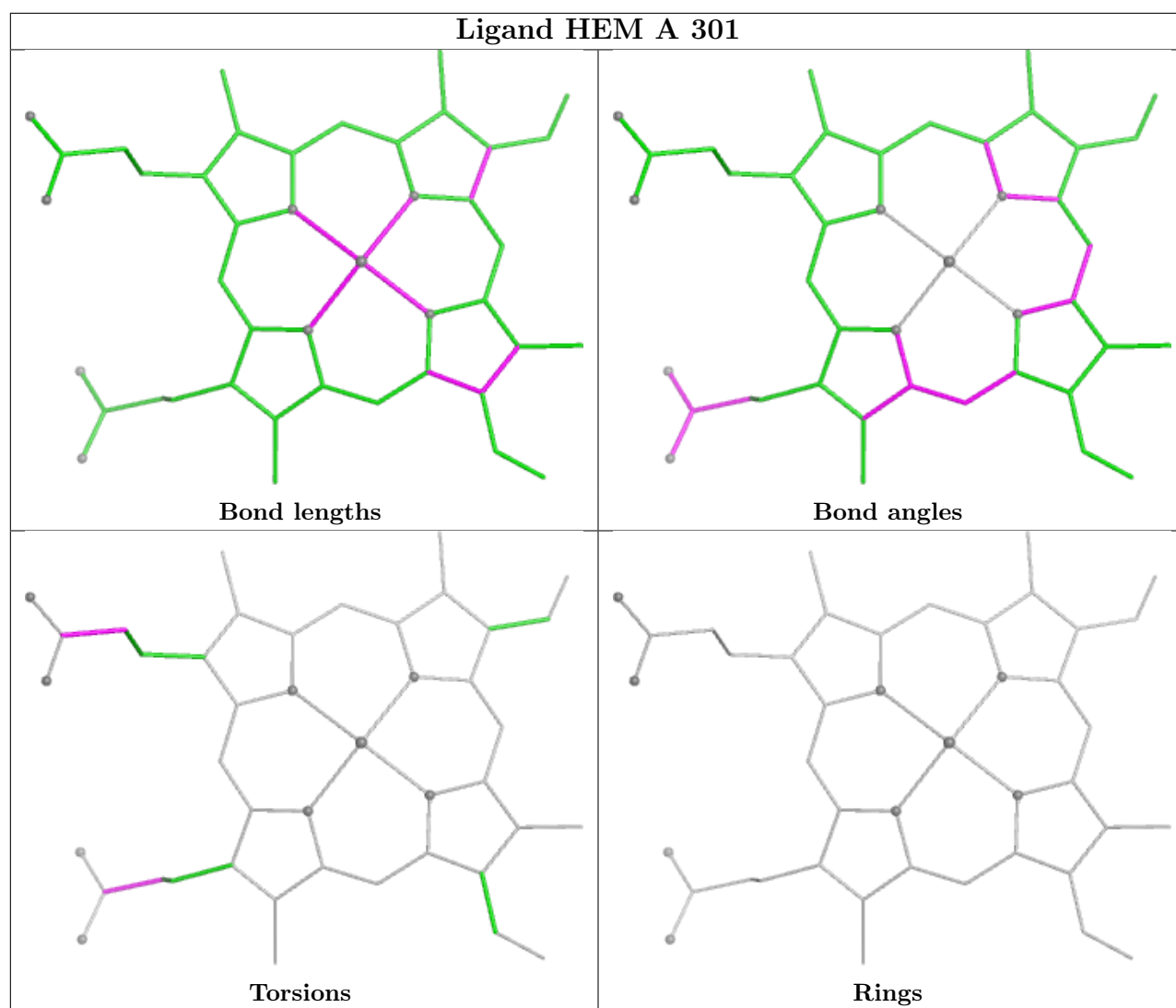
Mol	Chain	Res	Type	Atoms
7	B	803	NAG	O5-C5-C6-O6
7	B	803	NAG	C4-C5-C6-O6
5	A	302	PEG	O1-C1-C2-O2
5	A	302	PEG	O2-C3-C4-O4
7	B	802	NAG	C4-C5-C6-O6
4	A	301	HEM	CAD-CBD-CGD-O1D
4	A	301	HEM	CAD-CBD-CGD-O2D
4	A	301	HEM	CAA-CBA-CGA-O1A
4	A	301	HEM	CAA-CBA-CGA-O2A

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	302	PEG	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	100/114 (87%)	-0.81	2 (2%) 65 68	5, 12, 30, 44	1 (1%)
2	B	458/467 (98%)	-1.01	0 100 100	5, 11, 26, 44	6 (1%)
3	C	69/73 (94%)	-0.62	0 100 100	8, 15, 34, 47	2 (2%)
All	All	627/654 (95%)	-0.94	2 (0%) 90 92	5, 11, 27, 47	9 (1%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	271	ALA	2.6
1	A	239	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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	YCM	B	319	10/11	0.89	0.13	18,35,51,53	0
2	CSO	B	316	7/8	0.99	0.04	8,10,16,28	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

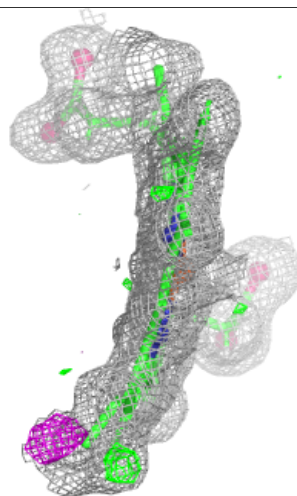
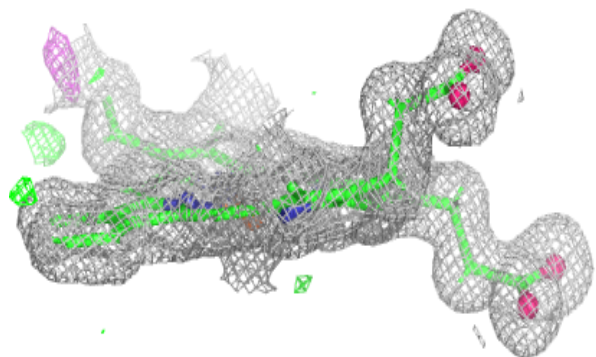
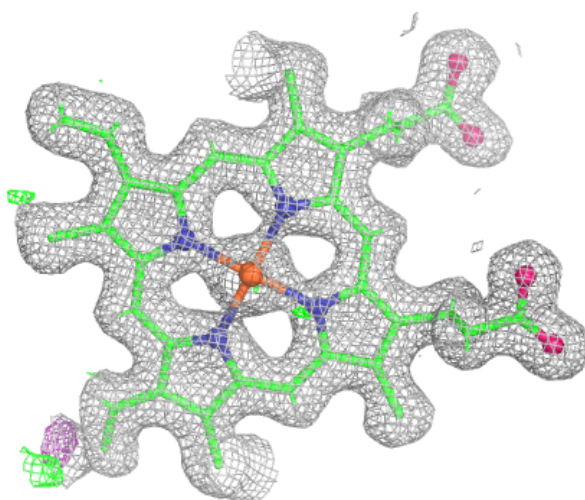
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	PEG	A	302	7/7	0.94	0.10	24,38,49,52	0
7	NAG	B	803	14/15	0.95	0.07	22,30,34,39	0
7	NAG	B	802	14/15	0.96	0.07	19,27,35,40	0
6	CL	B	810	1/1	0.98	0.09	47,47,47,47	0
7	NAG	B	801	14/15	0.99	0.03	7,11,15,17	0
9	MG	B	813	1/1	0.99	0.05	30,30,30,30	0
6	CL	B	808	1/1	1.00	0.04	19,19,19,19	0
6	CL	B	809	1/1	1.00	0.04	16,16,16,16	0
4	HEM	A	301	43/43	1.00	0.03	5,6,7,9	0
6	CL	B	811	1/1	1.00	0.13	37,37,37,37	0
6	CL	B	812	1/1	1.00	0.07	16,16,16,16	0
6	CL	A	303	1/1	1.00	0.01	8,8,8,8	0
6	CL	B	805	1/1	1.00	0.02	19,19,19,19	0
6	CL	B	806	1/1	1.00	0.03	21,21,21,21	0
8	CA	B	804	1/1	1.00	0.01	7,7,7,7	0
6	CL	B	807	1/1	1.00	0.02	21,21,21,21	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around HEM A 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.