



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 3, 2024 – 08:46 pm GMT

PDB ID : 4Z0Y
Title : Active aurone synthase (polyphenol oxidase), copper B : sulfohistidine 1.4 :
1
Authors : Molitor, C.; Mauracher, S.G.; Rompel, A.
Deposited on : 2015-03-26
Resolution : 1.60 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

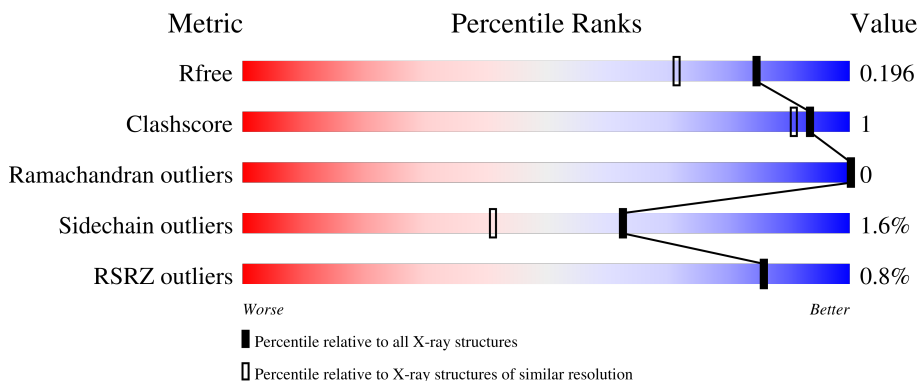
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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.



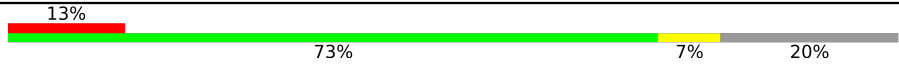

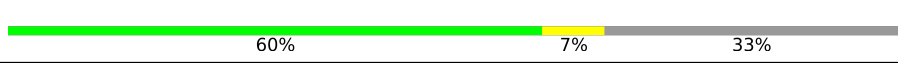
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	351	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 96%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">96%</p>
1	B	351	<div style="display: flex; align-items: center;"> <div style="width: 95%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">95%</p>
1	C	351	<div style="display: flex; align-items: center;"> <div style="width: 95%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">95%</p>
1	D	351	<div style="display: flex; align-items: center;"> <div style="width: 96%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">96%</p>
2	E	15	<div style="display: flex; align-items: center;"> <div style="width: 20%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 33%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 47%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">20% 53% 47%</p>

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Mol	Chain	Length	Quality of chain
2	F	15	 13% 73% 7% 20%
2	G	15	 13% 60% 7% 33%
2	H	15	 60% 7% 33%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 23988 atoms, of which 10376 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 Aurone synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	347	5317	1781	2513	478	532	13	0	4	0
1	B	347	5336	1777	2540	478	528	13	0	3	0
1	C	347	5381	1783	2574	480	531	13	0	4	0
1	D	347	5396	1790	2575	482	536	13	0	6	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	252	HS8	HIS	microheterogeneity	UNP A0A075DN54
B	252	HS8	HIS	microheterogeneity	UNP A0A075DN54
C	252	HS8	HIS	microheterogeneity	UNP A0A075DN54
D	252	HS8	HIS	microheterogeneity	UNP A0A075DN54

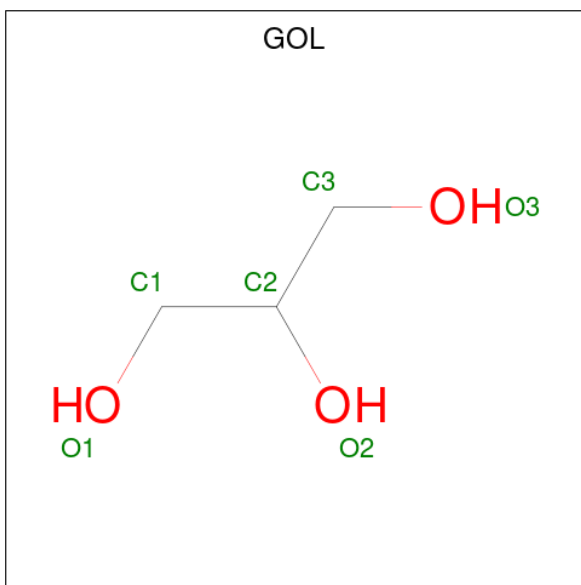
- Molecule 2 is a protein called Aurone synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	E	8	87	36	32	8	10	1	0	0	0
2	F	12	120	56	34	12	17	1	0	0	0
2	G	10	103	43	34	10	15	1	0	0	0
2	H	10	103	44	34	10	14	1	0	0	0

- Molecule 3 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cu 2 2	0	0
3	B	2	Total Cu 2 2	0	0
3	C	2	Total Cu 2 2	0	0
3	D	2	Total Cu 2 2	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C H O 14 3 8 3	0	0
4	B	1	Total C H O 14 3 8 3	0	0
4	C	1	Total C H O 14 3 8 3	0	0
4	D	1	Total C H O 14 3 8 3	0	0
4	D	1	Total C H O 14 3 8 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	463	Total O 463 463	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	9	Total 9	O 9	0	0
5	B	469	Total 469	O 469	0	0
5	F	21	Total 21	O 21	0	0
5	C	525	Total 525	O 525	0	0
5	G	11	Total 11	O 11	0	0
5	D	555	Total 555	O 555	0	0
5	H	14	Total 14	O 14	0	0

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: Aurone synthase



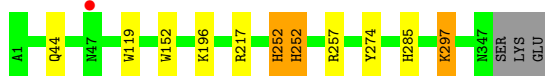
- Molecule 1: Aurone synthase



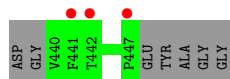
- Molecule 1: Aurone synthase



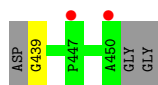
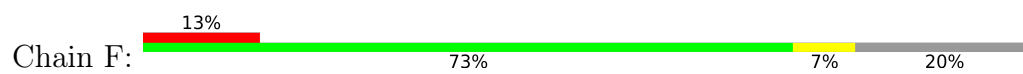
- Molecule 1: Aurone synthase



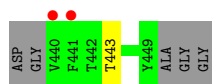
- Molecule 2: Aurone synthase



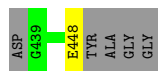
- Molecule 2: Aurone synthase



- Molecule 2: Aurone synthase



- Molecule 2: Aurone synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.72Å 90.56Å 182.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.65 – 1.60 45.65 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.3 (45.65-1.60) 99.4 (45.65-1.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 1.60Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.157 , 0.194 0.159 , 0.196	Depositor DCC
R_{free} test set	9628 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	14.9	Xtrriage
Anisotropy	0.589	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 51.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.000 for k,h,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	23988	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.34 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.5610e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CU, HS8

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	0.37	0/2875	0.55	0/3917
1	B	0.40	0/2867	0.57	0/3907
1	C	0.41	0/2878	0.58	0/3921
1	D	0.42	0/2892	0.59	0/3940
2	E	0.36	0/57	0.52	0/79
2	F	0.36	0/89	0.53	0/123
2	G	0.42	0/71	0.53	0/99
2	H	0.37	0/71	0.49	0/98
All	All	0.40	0/11800	0.57	0/16084

There are no bond length outliers.

There are no bond angle outliers.

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	2804	2513	2589	5	0
1	B	2796	2540	2583	10	0
1	C	2807	2574	2608	8	0
1	D	2821	2575	2611	7	0
2	E	55	32	44	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	86	34	72	1	0
2	G	69	34	50	0	0
2	H	69	34	58	1	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	6	8	8	0	0
4	B	6	8	8	3	0
4	C	6	8	8	0	0
4	D	12	16	16	2	0
5	A	463	0	0	3	1
5	B	469	0	0	4	1
5	C	525	0	0	5	2
5	D	555	0	0	4	1
5	E	9	0	0	0	1
5	F	21	0	0	1	0
5	G	11	0	0	0	0
5	H	14	0	0	1	0
All	All	13612	10376	10655	32	3

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 1.

All (32) 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:C:252[A]:HS8:O1	5:C:701:HOH:O	1.94	0.85
1:A:252[A]:HS8:O1	5:A:701:HOH:O	1.99	0.77
1:B:44:GLN:H	4:B:603:GOL:H12	1.51	0.74
1:A:252[A]:HS8:O2	5:A:702:HOH:O	2.06	0.74
1:B:112:ASP:OD2	5:B:703:HOH:O	2.07	0.73
1:B:252[A]:HS8:O1	5:B:702:HOH:O	2.05	0.72
1:B:44:GLN:H	4:B:603:GOL:C1	2.06	0.68
1:A:93:HIS:CE1	5:A:701:HOH:O	2.46	0.68
1:A:58:HIS:CD2	1:A:59:LYS:HG3	2.32	0.65
1:D:44:GLN:H	4:D:604:GOL:H32	1.63	0.64
1:D:44:GLN:H	4:D:604:GOL:C3	2.11	0.63
2:F:439:GLY:N	5:F:502:HOH:O	2.33	0.62
1:D:196:LYS:O	5:D:702:HOH:O	2.16	0.60
1:D:297:LYS:NZ	5:D:706:HOH:O	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:311:ASP:OD2	5:C:703:HOH:O	2.18	0.56
1:C:70:ARG:NH2	5:C:709:HOH:O	2.40	0.54
1:C:252[A]:HS8:O2	5:C:702:HOH:O	2.17	0.54
1:C:252[A]:HS8:S	5:C:701:HOH:O	2.52	0.54
1:D:252[A]:HS8:HD1	1:D:285:HIS:HE2	1.57	0.52
1:B:84:ARG:NH1	5:B:716:HOH:O	2.41	0.51
1:D:252[A]:HS8:O2	5:D:703:HOH:O	2.20	0.48
1:B:325:ASP:OD1	5:B:704:HOH:O	2.20	0.48
1:B:221:ASP:HB2	1:B:222:PRO:CD	2.47	0.44
1:C:58:HIS:CD2	1:C:59:LYS:HG3	2.53	0.44
1:C:296:TRP:CH2	1:C:304:HIS:HA	2.52	0.44
1:C:252[A]:HS8:HD1	1:C:285:HIS:HE2	1.64	0.44
2:H:448:GLU:HA	5:H:511:HOH:O	2.17	0.44
1:D:257:ARG:NH1	5:D:709:HOH:O	2.37	0.43
1:B:127:TRP:CE2	1:B:334:CYS:HB3	2.54	0.42
1:A:296:TRP:CH2	1:A:304:HIS:HA	2.53	0.42
1:B:221:ASP:HB2	1:B:222:PRO:HD2	2.03	0.41
1:B:44:GLN:N	4:B:603:GOL:H12	2.29	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:1066:HOH:O	5:C:928:HOH:O[2_575]	2.09	0.11
5:A:1088:HOH:O	5:D:1060:HOH:O[4_477]	2.13	0.07
5:E:502:HOH:O	5:C:806:HOH:O[3_557]	2.14	0.06

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	347/351 (99%)	335 (96%)	12 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	346/351 (99%)	338 (98%)	8 (2%)	0	100	100
1	C	347/351 (99%)	338 (97%)	9 (3%)	0	100	100
1	D	349/351 (99%)	339 (97%)	10 (3%)	0	100	100
2	E	6/15 (40%)	6 (100%)	0	0	100	100
2	F	10/15 (67%)	10 (100%)	0	0	100	100
2	G	8/15 (53%)	7 (88%)	1 (12%)	0	100	100
2	H	8/15 (53%)	8 (100%)	0	0	100	100
All	All	1421/1464 (97%)	1381 (97%)	40 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	294/305 (96%)	291 (99%)	3 (1%)	76	61
1	B	293/305 (96%)	289 (99%)	4 (1%)	67	47
1	C	297/305 (97%)	291 (98%)	6 (2%)	55	31
1	D	298/305 (98%)	293 (98%)	5 (2%)	60	38
2	E	6/11 (54%)	6 (100%)	0	100	100
2	F	9/11 (82%)	9 (100%)	0	100	100
2	G	7/11 (64%)	6 (86%)	1 (14%)	3	0
2	H	8/11 (73%)	8 (100%)	0	100	100
All	All	1212/1264 (96%)	1193 (98%)	19 (2%)	62	41

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	TRP
1	A	152	TRP

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Mol	Chain	Res	Type
1	A	274	TYR
1	B	48	ASP
1	B	119	TRP
1	B	152	TRP
1	B	274	TYR
1	C	48	ASP
1	C	119	TRP
1	C	152	TRP
1	C	274	TYR
1	C	281	VAL
1	C	317	TYR
2	G	443	THR
1	D	119	TRP
1	D	152	TRP
1	D	217	ARG
1	D	274	TYR
1	D	297	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	44	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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	HS8	D	252[A]	3,1	10,14,15	2.16	4 (40%)	4,20,22	2.79	2 (50%)
1	HS8	B	252[A]	3,1	10,14,15	4.08	3 (30%)	4,20,22	3.02	2 (50%)
1	HS8	A	252[A]	3,1	10,14,15	2.24	4 (40%)	4,20,22	2.32	2 (50%)
1	HS8	C	252[A]	3,1	10,14,15	2.23	4 (40%)	4,20,22	2.52	2 (50%)

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	HS8	D	252[A]	3,1	-	0/5/12/14	0/1/1/1
1	HS8	B	252[A]	3,1	-	0/5/12/14	0/1/1/1
1	HS8	A	252[A]	3,1	-	0/5/12/14	0/1/1/1
1	HS8	C	252[A]	3,1	-	0/5/12/14	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	252[A]	HS8	O1-S	11.82	1.62	1.42
1	A	252[A]	HS8	O1-S	-3.85	1.36	1.42
1	D	252[A]	HS8	O1-S	-3.82	1.36	1.42
1	C	252[A]	HS8	O1-S	-3.70	1.36	1.42
1	C	252[A]	HS8	O3-S	-3.52	1.36	1.42
1	A	252[A]	HS8	O2-S	3.31	1.62	1.45
1	D	252[A]	HS8	O2-S	3.30	1.62	1.45
1	C	252[A]	HS8	O2-S	3.30	1.62	1.45
1	A	252[A]	HS8	O3-S	-3.20	1.37	1.42
1	D	252[A]	HS8	O3-S	-3.17	1.37	1.42
1	B	252[A]	HS8	CD2-CG	3.09	1.40	1.36
1	B	252[A]	HS8	O3-S	-2.94	1.37	1.42
1	A	252[A]	HS8	CD2-CG	2.84	1.40	1.36
1	C	252[A]	HS8	CD2-CG	2.80	1.40	1.36
1	D	252[A]	HS8	CD2-CG	2.58	1.40	1.36

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	252[A]	HS8	O3-S-O1	-5.35	105.29	119.22
1	C	252[A]	HS8	O3-S-O1	4.36	130.58	119.22
1	D	252[A]	HS8	O3-S-O1	4.30	130.41	119.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	252[A]	HS8	O3-S-O1	3.54	128.42	119.22
1	D	252[A]	HS8	CB-CA-C	-3.26	105.36	111.47
1	A	252[A]	HS8	CB-CA-C	-2.66	106.48	111.47
1	B	252[A]	HS8	CB-CA-C	-2.60	106.59	111.47
1	C	252[A]	HS8	CB-CA-C	-2.34	107.08	111.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	252[A]	HS8	2	0
1	B	252[A]	HS8	1	0
1	A	252[A]	HS8	2	0
1	C	252[A]	HS8	4	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 8 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	$\# Z > 2$
4	GOL	B	603	-	5,5,5	0.28	0	5,5,5	0.48	0
4	GOL	D	603	-	5,5,5	0.29	0	5,5,5	0.55	0
4	GOL	A	603	-	5,5,5	0.28	0	5,5,5	0.28	0
4	GOL	C	603	-	5,5,5	0.41	0	5,5,5	0.71	0
4	GOL	D	604	-	5,5,5	0.24	0	5,5,5	0.48	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	GOL	B	603	-	-	2/4/4/4	-
4	GOL	D	603	-	-	1/4/4/4	-
4	GOL	A	603	-	-	0/4/4/4	-
4	GOL	C	603	-	-	2/4/4/4	-
4	GOL	D	604	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	603	GOL	O1-C1-C2-C3
4	B	603	GOL	O1-C1-C2-O2
4	C	603	GOL	O1-C1-C2-C3
4	C	603	GOL	O1-C1-C2-O2
4	D	603	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	603	GOL	3	0
4	D	604	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	346/351 (98%)	-0.47	2 (0%) 89 89	11, 19, 34, 63	0
1	B	346/351 (98%)	-0.48	1 (0%) 94 93	9, 16, 30, 47	0
1	C	346/351 (98%)	-0.52	0 100 100	10, 17, 30, 48	0
1	D	346/351 (98%)	-0.60	1 (0%) 94 93	8, 13, 25, 52	0
2	E	8/15 (53%)	1.72	3 (37%) 0 0	34, 44, 52, 73	0
2	F	12/15 (80%)	0.83	2 (16%) 1 1	17, 26, 72, 86	0
2	G	10/15 (66%)	1.35	2 (20%) 1 1	26, 38, 62, 80	0
2	H	10/15 (66%)	-0.29	0 100 100	13, 17, 43, 57	0
All	All	1424/1464 (97%)	-0.48	11 (0%) 86 86	8, 16, 32, 86	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	441	PHE	5.0
2	F	450	ALA	4.6
1	A	198	ILE	4.3
2	G	440	VAL	3.6
2	E	442	THR	3.0
2	E	441	PHE	2.6
2	F	447	PRO	2.6
1	A	212	TYR	2.5
1	B	47	ASN	2.4
2	E	447	PRO	2.4
1	D	47	ASN	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
1	HS8	A	252[A]	14/15	0.93	0.11	14,16,28,36	14
1	HS8	B	252[A]	14/15	0.94	0.10	10,13,27,32	14
1	HS8	C	252[A]	14/15	0.97	0.09	11,14,23,36	14
1	HS8	D	252[A]	14/15	0.97	0.09	9,11,25,30	14

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
4	GOL	B	603	6/6	0.77	0.25	21,34,46,48	0
4	GOL	C	603	6/6	0.80	0.18	25,34,41,48	0
4	GOL	D	604	6/6	0.81	0.13	24,36,43,44	0
4	GOL	A	603	6/6	0.88	0.12	21,28,34,34	0
4	GOL	D	603	6/6	0.92	0.11	17,23,28,28	0
3	CU	B	602	1/1	0.96	0.04	23,23,23,23	1
3	CU	C	602	1/1	0.96	0.05	23,23,23,23	1
3	CU	D	602	1/1	0.96	0.05	22,22,22,22	1
3	CU	A	602	1/1	0.96	0.04	24,24,24,24	1
3	CU	A	601	1/1	0.99	0.03	18,18,18,18	0
3	CU	B	601	1/1	0.99	0.04	17,17,17,17	0
3	CU	C	601	1/1	1.00	0.04	16,16,16,16	0
3	CU	D	601	1/1	1.00	0.04	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.