



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:03 AM UTC

PDB ID : 4Z0Z / pdb_00004z0z
Title : Inactive aurone synthase (polyphenol oxidase) from natural source, sulfohistidine 90 %
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:

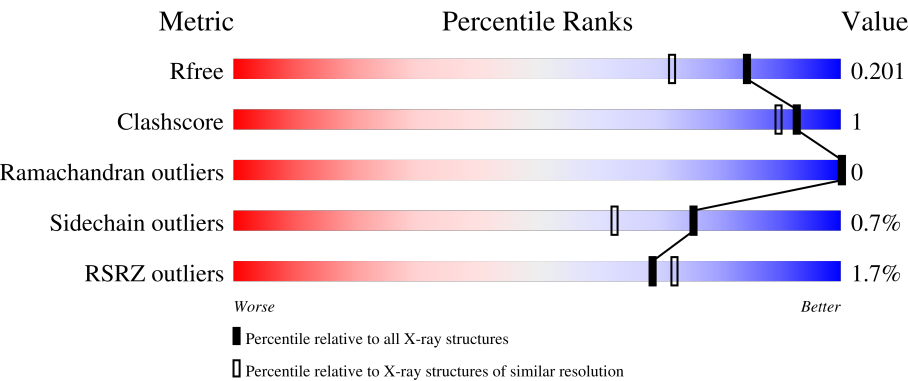
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (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	350	96% ..
1	B	350	% 96% ..
1	C	350	% 95% ..
1	D	350	% 94% 5%
2	E	15	27% 60% 7% 33%

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Mol	Chain	Length	Quality of chain
2	F	15	
2	G	15	
2	H	15	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23779 atoms, of which 10327 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
1	A	347	Total	C	H	N	O	S	0	0	0
			5344	1772	2555	476	528	13			
1	B	347	Total	C	H	N	O	S	0	1	0
			5332	1770	2550	474	525	13			
1	C	347	Total	C	H	N	O	S	0	0	0
			5327	1769	2547	475	523	13			
1	D	347	Total	C	H	N	O	S	0	1	0
			5331	1771	2547	474	526	13			

- Molecule 2 is a protein called Aurone synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	10	Total	C	H	N	O	S	0	0	0
			111	51	34	10	15	1			
2	F	9	Total	C	H	N	O	S	0	0	0
			100	40	36	9	14	1			
2	G	9	Total	C	H	N	O	S	0	0	0
			87	42	22	9	13	1			
2	H	9	Total	C	H	N	O	S	0	0	0
			100	41	36	9	13	1			

- Molecule 3 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

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

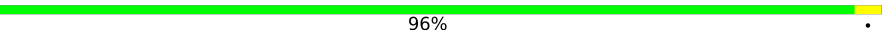
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	485	Total 485	O 485	0	0
4	E	13	Total 13	O 13	0	0
4	B	518	Total 518	O 518	0	0
4	F	16	Total 16	O 16	0	0
4	C	469	Total 469	O 469	0	0
4	G	13	Total 13	O 13	0	0
4	D	511	Total 511	O 511	0	0
4	H	18	Total 18	O 18	0	0

3 Residue-property plots [i](#)

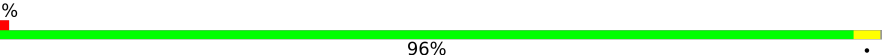
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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

Chain A:  96% ..

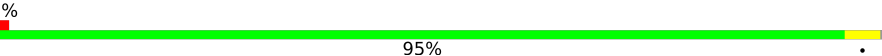


- Molecule 1: Aurone synthase

Chain B:  96% ..

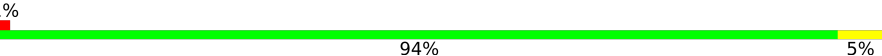


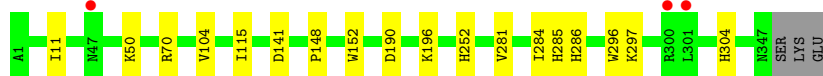
- Molecule 1: Aurone synthase

Chain C:  95% ..



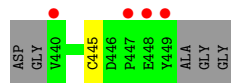
- Molecule 1: Aurone synthase

Chain D:  94% 5% .

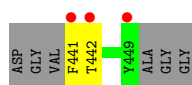


- Molecule 2: Aurone synthase

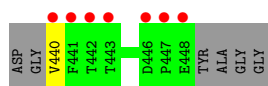
Chain E:  27% 60% 7% 33%



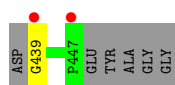
- Molecule 2: Aurone synthase



- Molecule 2: Aurone synthase



- Molecule 2: Aurone synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.53Å 183.52Å 78.09Å 90.00° 94.50° 90.00°	Depositor
Resolution (Å)	48.10 – 1.60 48.10 – 1.60	Depositor EDS
% Data completeness (in resolution range)	98.7 (48.10-1.60) 98.7 (48.10-1.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 1.60Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.165 , 0.200 0.167 , 0.201	Depositor DCC
R_{free} test set	9354 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	16.2	Xtriage
Anisotropy	0.395	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	23779	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.56% 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: 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	0.46	0/2859	0.67	2/3894 (0.1%)
1	B	0.47	0/2855	0.69	0/3889
1	C	0.46	0/2850	0.69	0/3882
1	D	0.49	0/2857	0.70	1/3892 (0.0%)
2	E	0.38	0/80	0.61	0/111
2	F	0.38	0/66	0.68	0/92
2	G	0.33	0/67	0.73	0/93
2	H	0.35	0/66	0.65	0/91
All	All	0.47	0/11700	0.69	3/15944 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	196	LYS	N-CA-C	5.31	117.98	111.82
1	A	122	PHE	CA-C-N	-5.03	113.86	119.19
1	A	122	PHE	C-N-CA	-5.03	113.86	119.19

There are no chirality outliers.

There are no planarity outliers.

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	2789	2555	2598	7	0
1	B	2782	2550	2591	6	0
1	C	2780	2547	2590	7	0
1	D	2784	2547	2589	9	0
2	E	77	34	64	1	0
2	F	64	36	48	1	0
2	G	65	22	55	1	0
2	H	64	36	56	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	485	0	0	4	0
4	B	518	0	0	1	3
4	C	469	0	0	4	4
4	D	511	0	0	4	4
4	E	13	0	0	1	0
4	F	16	0	0	1	0
4	G	13	0	0	0	0
4	H	18	0	0	1	0
All	All	13452	10327	10591	32	6

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:112:ASP:OD2	4:C:801:HOH:O	2.03	0.73
1:D:141:ASP:OD2	4:D:801:HOH:O	2.06	0.72
1:C:195:ASP:OD2	4:C:802:HOH:O	2.12	0.67
2:H:439:GLY:N	4:H:502:HOH:O	2.26	0.66
2:E:445:CYS:HB2	4:E:505:HOH:O	1.96	0.65
1:D:70:ARG:NH1	4:D:805:HOH:O	2.26	0.65
2:F:441:PHE:N	4:F:501:HOH:O	2.29	0.64
1:D:152:TRP:HE1	1:D:285:HIS:HD1	1.49	0.61
1:C:152:TRP:HE1	1:C:285:HIS:HD1	1.47	0.60
1:C:311:ASP:OD1	4:C:803:HOH:O	2.16	0.58
1:A:156:LYS:NZ	4:A:809:HOH:O	2.34	0.57
1:D:50:LYS:HG3	4:D:1106:HOH:O	2.05	0.56
1:B:299:LEU:HD23	2:G:440:VAL:HG21	1.90	0.54
1:C:297:LYS:HE3	1:C:306:ASP:OD2	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:TRP:HE1	1:B:285:HIS:HD1	1.56	0.52
1:D:190:ASP:OD1	4:D:802:HOH:O	2.19	0.49
1:A:105:ASP:OD1	4:A:801:HOH:O	2.20	0.47
1:A:21:ASN:OD1	4:A:802:HOH:O	2.20	0.47
1:C:221:ASP:HB2	1:C:222:PRO:CD	2.46	0.46
1:A:44:GLN:OE1	4:A:803:HOH:O	2.21	0.46
1:A:252:HS8:HD1	1:A:285:HIS:HE2	1.63	0.45
1:A:11:ILE:HD13	1:A:104:VAL:HG11	1.98	0.45
1:D:11:ILE:HD13	1:D:104:VAL:HG11	1.99	0.45
1:C:198:ILE:O	4:C:804:HOH:O	2.20	0.45
1:D:148:PRO:HG2	1:D:284:ILE:HD12	1.99	0.43
1:B:125:HIS:CG	1:B:286:HIS:CE1	3.07	0.43
1:A:174:ASP:OD2	1:A:269:ASP:OD2	2.36	0.42
1:D:115:ILE:C	1:D:115:ILE:HD12	2.45	0.42
1:D:296:TRP:CH2	1:D:304:HIS:HA	2.55	0.42
1:B:190:ASP:OD1	4:B:801:HOH:O	2.21	0.41
1:B:296:TRP:CH2	1:B:304:HIS:HA	2.56	0.41
1:B:221:ASP:HB2	1:B:222:PRO:CD	2.51	0.41

All (6) 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)
4:C:1051:HOH:O	4:C:1111:HOH:O[1_455]	2.10	0.10
4:B:1120:HOH:O	4:D:1224:HOH:O[1_556]	2.13	0.07
4:C:870:HOH:O	4:D:890:HOH:O[2_545]	2.13	0.07
4:B:1312:HOH:O	4:C:1258:HOH:O[2_555]	2.14	0.06
4:C:1131:HOH:O	4:D:890:HOH:O[2_545]	2.14	0.06
4:B:1212:HOH:O	4:D:980:HOH:O[1_556]	2.19	0.01

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	344/350 (98%)	333 (97%)	11 (3%)	0	100	100
1	B	345/350 (99%)	334 (97%)	11 (3%)	0	100	100
1	C	344/350 (98%)	333 (97%)	11 (3%)	0	100	100
1	D	345/350 (99%)	333 (96%)	12 (4%)	0	100	100
2	E	8/15 (53%)	7 (88%)	1 (12%)	0	100	100
2	F	7/15 (47%)	7 (100%)	0	0	100	100
2	G	7/15 (47%)	7 (100%)	0	0	100	100
2	H	7/15 (47%)	7 (100%)	0	0	100	100
All	All	1407/1460 (96%)	1361 (97%)	46 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	295/304 (97%)	294 (100%)	1 (0%)	86	78
1	B	293/304 (96%)	291 (99%)	2 (1%)	76	63
1	C	292/304 (96%)	290 (99%)	2 (1%)	76	63
1	D	293/304 (96%)	290 (99%)	3 (1%)	68	50
2	E	9/11 (82%)	9 (100%)	0	100	100
2	F	7/11 (64%)	6 (86%)	1 (14%)	3	0
2	G	8/11 (73%)	8 (100%)	0	100	100
2	H	8/11 (73%)	8 (100%)	0	100	100
All	All	1205/1260 (96%)	1196 (99%)	9 (1%)	76	63

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	286	HIS
1	B	281	VAL

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Mol	Chain	Res	Type
1	B	286	HIS
2	F	442	THR
1	C	36	LEU
1	C	286	HIS
1	D	281	VAL
1	D	286	HIS
1	D	297	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	323	ASN
1	B	98	ASN
1	B	117	ASN
1	B	144	ASN
1	B	202	GLN
1	B	239	ASN
1	B	264	GLN
1	B	314	ASN
1	C	264	GLN
1	D	98	ASN

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	HS8	A	252	1	12,14,15	2.09	4 (33%)	13,20,22	1.50	3 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HS8	C	252	1	12,14,15	2.16	4 (33%)	13,20,22	2.18	5 (38%)
1	HS8	B	252	1	12,14,15	1.97	3 (25%)	13,20,22	1.62	4 (30%)
1	HS8	D	252	1	12,14,15	4.91	4 (33%)	13,20,22	2.15	3 (23%)

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	A	252	1	-	3/9/12/14	0/1/1/1
1	HS8	C	252	1	-	0/9/12/14	0/1/1/1
1	HS8	B	252	1	-	3/9/12/14	0/1/1/1
1	HS8	D	252	1	-	0/9/12/14	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	252	HS8	O1-S	15.96	1.60	1.42
1	B	252	HS8	O3-S	-4.50	1.37	1.42
1	C	252	HS8	O3-S	-4.19	1.37	1.42
1	A	252	HS8	O1-S	-4.16	1.37	1.42
1	D	252	HS8	O3-S	-3.99	1.37	1.42
1	A	252	HS8	O3-S	-3.70	1.37	1.42
1	C	252	HS8	O1-S	-3.66	1.37	1.42
1	B	252	HS8	O1-S	-2.96	1.38	1.42
1	A	252	HS8	O2-S	2.96	1.61	1.47
1	C	252	HS8	O2-S	2.94	1.61	1.47
1	B	252	HS8	O2-S	2.83	1.61	1.47
1	C	252	HS8	CD2-CG	2.79	1.41	1.36
1	D	252	HS8	CD2-CG	2.63	1.40	1.36
1	A	252	HS8	CD2-CG	2.60	1.40	1.36
1	D	252	HS8	O2-S	-2.28	1.36	1.47

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	252	HS8	O3-S-O1	-5.08	107.35	119.69
1	D	252	HS8	O2-S-O3	4.39	127.48	113.10
1	C	252	HS8	CE1-NE2-S	-4.22	123.22	126.55
1	C	252	HS8	O3-S-O1	3.75	128.79	119.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	252	HS8	CD2-NE2-S	3.19	129.20	126.09
1	B	252	HS8	O3-S-O1	3.18	127.41	119.69
1	C	252	HS8	O2-S-O3	-2.80	103.91	113.10
1	A	252	HS8	O3-S-O1	2.40	125.53	119.69
1	A	252	HS8	O2-S-O3	-2.28	105.62	113.10
1	A	252	HS8	CE1-NE2-S	-2.22	124.79	126.55
1	D	252	HS8	CA-CB-CG	-2.21	108.31	113.77
1	B	252	HS8	CE1-ND1-CG	2.19	109.31	105.80
1	B	252	HS8	NE2-CE1-ND1	-2.04	108.47	111.68
1	C	252	HS8	CE1-ND1-CG	2.02	109.04	105.80
1	B	252	HS8	CD2-NE2-CE1	2.00	108.78	106.66

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	252	HS8	CE1-NE2-S-O1
1	B	252	HS8	CD2-NE2-S-O1
1	A	252	HS8	CD2-NE2-S-O3
1	A	252	HS8	CE1-NE2-S-O3
1	A	252	HS8	C-CA-CB-CG
1	B	252	HS8	CE1-NE2-S-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	252	HS8	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	346/350 (98%)	-0.29	0	100	100	11, 20, 35, 52	0
1	B	346/350 (98%)	-0.34	3 (0%)	81	84	9, 19, 35, 60	1 (0%)
1	C	346/350 (98%)	-0.29	2 (0%)	85	88	12, 20, 36, 61	0
1	D	346/350 (98%)	-0.48	3 (0%)	81	84	8, 17, 31, 64	1 (0%)
2	E	10/15 (66%)	1.61	4 (40%)	1	0	28, 34, 53, 70	0
2	F	9/15 (60%)	1.73	3 (33%)	1	1	28, 39, 52, 52	0
2	G	9/15 (60%)	2.54	7 (77%)	0	0	32, 36, 65, 89	0
2	H	9/15 (60%)	0.97	2 (22%)	2	2	20, 21, 57, 66	0
All	All	1421/1460 (97%)	-0.30	24 (1%)	69	72	8, 19, 36, 89	2 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	441	PHE	6.3
2	G	447	PRO	5.0
1	C	197	ASP	4.1
1	C	212	TYR	4.0
2	G	441	PHE	3.8
2	H	439	GLY	3.7
2	H	447	PRO	3.6
2	E	440	VAL	3.2
2	E	448	GLU	3.0
1	B	47	ASN	2.9
2	G	448	GLU	2.7
1	D	300	ARG	2.7
2	G	440	VAL	2.5
1	B	49	LYS	2.4
2	G	442	THR	2.4
1	B	212	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
2	G	446	ASP	2.4
2	E	447	PRO	2.3
1	D	301	LEU	2.3
2	F	442	THR	2.2
1	D	47	ASN	2.1
2	E	449	TYR	2.1
2	G	443	THR	2.0
2	F	449	TYR	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. 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	B	252	14/15	0.96	0.07	10,13,23,25	4
1	HS8	A	252	14/15	0.97	0.06	13,16,24,29	4
1	HS8	C	252	14/15	0.97	0.06	12,15,25,26	4
1	HS8	D	252	14/15	0.97	0.06	10,12,20,24	4

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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
3	CU	A	700	1/1	1.00	0.02	17,17,17,17	0
3	CU	B	700	1/1	1.00	0.01	15,15,15,15	0
3	CU	C	700	1/1	1.00	0.03	17,17,17,17	0
3	CU	D	700	1/1	1.00	0.02	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.