



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:49 PM GMT

PDB ID : 1HDH
Title : ARYLSULFATASE FROM PSEUDOMONAS AERUGINOSA
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Deposited on : 2000-11-16
Resolution : 1.30 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

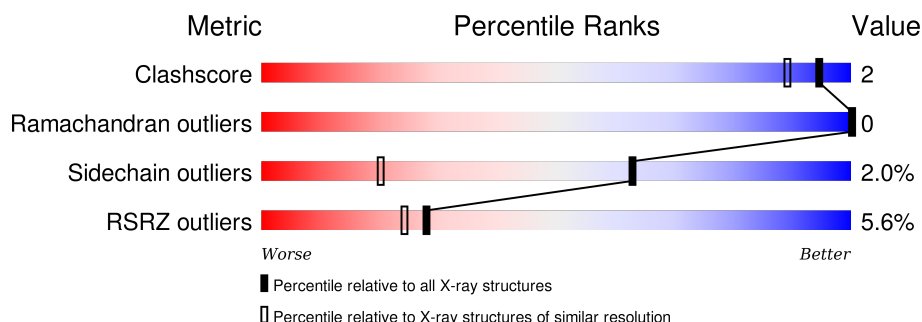
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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))
Clashscore	102246	1031 (1.32-1.28)
Ramachandran outliers	100387	1504 (1.34-1.26)
Sidechain outliers	100360	1503 (1.34-1.26)
RSRZ outliers	91569	1476 (1.34-1.26)

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	536	<div> <div>5%</div> <div>89%</div> <div>8%</div> <div>..</div> </div>
1	B	536	<div> <div>6%</div> <div>89%</div> <div>8%</div> <div>..</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
3	SO4	A	1529	-	-	-	X
3	SO4	B	1529	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	B	1532	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9041 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 ARYLSULFATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	525	Total	C	N	O	S	0	5	0
			4095	2613	720	753	9			
1	B	525	Total	C	N	O	S	0	7	0
			4094	2610	719	756	9			

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

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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

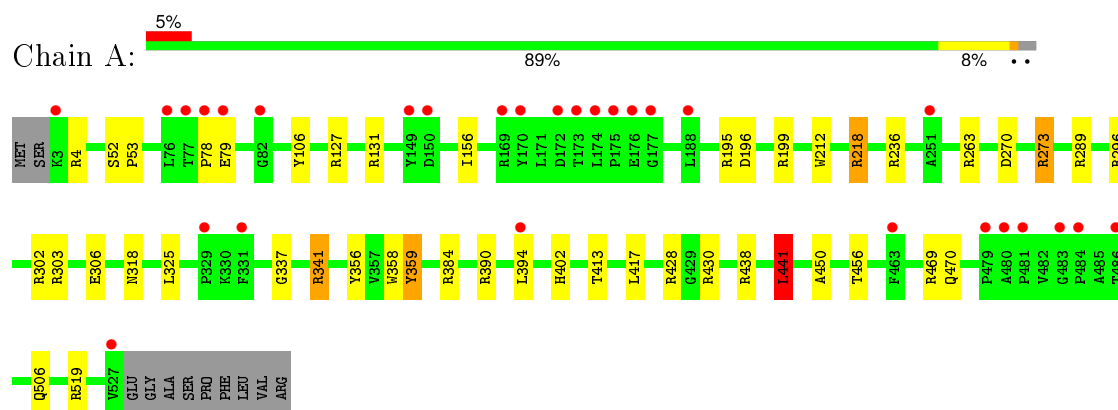
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	412	Total O 412 412	0	0
4	B	398	Total O 398 398	0	0

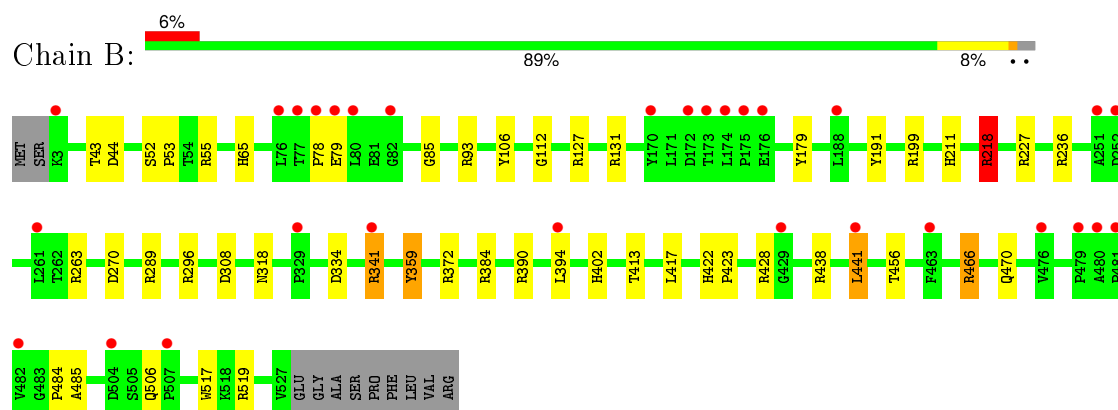
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: ARYLSULFATASE



• Molecule 1: ARYLSULFATASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	188.92Å 67.29Å 89.70Å 90.00° 94.20° 90.00°	Depositor
Resolution (Å)	20.00 – 1.30 19.99 – 1.30	Depositor EDS
% Data completeness (in resolution range)	93.1 (20.00-1.30) 93.1 (19.99-1.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 1.30Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.200 , 0.229 0.207 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	10.4	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 255817 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9041	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DDZ, CA, SO4

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	0.57	0/4223	1.27	37/5751 (0.6%)
1	B	0.57	0/4233	1.26	37/5767 (0.6%)
All	All	0.57	0/8456	1.27	74/11518 (0.6%)

There are no bond length outliers.

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	236	ARG	NE-CZ-NH1	16.37	128.48	120.30
1	A	273	ARG	CD-NE-CZ	11.75	140.04	123.60
1	A	263	ARG	NE-CZ-NH1	11.66	126.13	120.30
1	B	390	ARG	NE-CZ-NH1	11.40	126.00	120.30
1	A	390	ARG	NE-CZ-NH1	11.06	125.83	120.30
1	A	390	ARG	NE-CZ-NH2	-9.87	115.37	120.30
1	B	227	ARG	NE-CZ-NH2	-9.25	115.67	120.30
1	B	236	ARG	NE-CZ-NH1	8.96	124.78	120.30
1	A	236	ARG	NE-CZ-NH2	-8.75	115.93	120.30
1	B	341[A]	ARG	CD-NE-CZ	8.69	135.77	123.60
1	B	341[B]	ARG	CD-NE-CZ	8.69	135.77	123.60
1	B	218	ARG	NE-CZ-NH2	8.64	124.62	120.30
1	A	263	ARG	NE-CZ-NH2	-8.59	116.01	120.30
1	B	263	ARG	CD-NE-CZ	8.57	135.60	123.60
1	A	341[A]	ARG	CD-NE-CZ	8.56	135.58	123.60
1	A	341[B]	ARG	CD-NE-CZ	8.56	135.58	123.60
1	A	199	ARG	NE-CZ-NH1	8.35	124.48	120.30
1	B	199	ARG	NE-CZ-NH1	8.15	124.38	120.30
1	A	430	ARG	NE-CZ-NH1	7.97	124.28	120.30
1	A	127	ARG	CD-NE-CZ	7.78	134.49	123.60
1	A	438[A]	ARG	NE-CZ-NH1	7.68	124.14	120.30
1	A	438[B]	ARG	NE-CZ-NH1	7.68	124.14	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	359	TYR	CB-CG-CD1	7.67	125.60	121.00
1	B	428	ARG	NE-CZ-NH1	7.24	123.92	120.30
1	B	296	ARG	NE-CZ-NH2	-7.23	116.68	120.30
1	A	428	ARG	NE-CZ-NH1	7.09	123.84	120.30
1	B	438	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	A	359	TYR	CB-CG-CD1	6.64	124.98	121.00
1	B	106	TYR	CZ-CE2-CD2	6.38	125.54	119.80
1	A	302	ARG	NE-CZ-NH2	-6.25	117.17	120.30
1	B	106	TYR	CG-CD1-CE1	6.24	126.29	121.30
1	B	179	TYR	CB-CG-CD1	-6.24	117.26	121.00
1	B	341[A]	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	B	341[B]	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	A	303	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	B	131	ARG	NE-CZ-NH2	-6.13	117.23	120.30
1	B	384	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	B	236	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	B	466	ARG	NE-CZ-NH2	-6.05	117.27	120.30
1	B	191	TYR	CB-CG-CD1	6.02	124.61	121.00
1	B	55	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	B	93	ARG	NE-CZ-NH1	5.96	123.28	120.30
1	B	127	ARG	CD-NE-CZ	5.93	131.91	123.60
1	A	106	TYR	CB-CG-CD1	5.93	124.56	121.00
1	A	4	ARG	NE-CZ-NH1	-5.92	117.34	120.30
1	A	199	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	A	384	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	A	106	TYR	CG-CD1-CE1	5.83	125.96	121.30
1	A	306	GLU	OE1-CD-OE2	-5.67	116.49	123.30
1	A	263	ARG	CD-NE-CZ	5.66	131.53	123.60
1	A	441	LEU	CB-CG-CD2	5.64	120.59	111.00
1	B	289	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	A	196	ASP	CB-CG-OD1	5.55	123.30	118.30
1	A	519	ARG	NE-CZ-NH1	5.55	123.07	120.30
1	B	44	ASP	CB-CG-OD1	5.50	123.25	118.30
1	B	390	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	B	218	ARG	NH1-CZ-NH2	-5.39	113.47	119.40
1	A	359	TYR	CA-CB-CG	5.34	123.55	113.40
1	B	296	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	A	469	ARG	CD-NE-CZ	5.22	130.91	123.60
1	A	519	ARG	CD-NE-CZ	5.22	130.90	123.60
1	A	273	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	B	308	ASP	CB-CG-OD1	5.16	122.94	118.30
1	A	296	ARG	NE-CZ-NH1	5.12	122.86	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	211	HIS	CA-CB-CG	5.11	122.28	113.60
1	B	131	ARG	NH1-CZ-NH2	5.09	125.00	119.40
1	B	372	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	A	289	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	A	356	TYR	CB-CG-CD2	5.08	124.05	121.00
1	B	131	ARG	NE-CZ-NH1	-5.08	117.76	120.30
1	B	263	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	A	131	ARG	NE-CZ-NH1	-5.06	117.77	120.30
1	A	195	ARG	NE-CZ-NH1	5.05	122.82	120.30
1	B	55	ARG	NE-CZ-NH2	-5.02	117.79	120.30

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	4095	0	3911	13	0
1	B	4094	0	3896	14	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	20	0	0	0	0
3	B	20	0	0	0	0
4	A	412	0	0	2	0
4	B	398	0	0	0	0
All	All	9041	0	7807	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 2.

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)
1:A:417:LEU:HG	1:A:441:LEU:HD21	1.72	0.71
1:A:337:GLY:O	1:A:341[B]:ARG:HD3	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:LEU:HG	1:B:441:LEU:HD21	1.82	0.62
1:A:78:PRO:HG2	1:A:79:GLU:OE2	2.03	0.57
1:B:78:PRO:HG2	1:B:79:GLU:OE2	2.05	0.56
1:B:413:THR:HG23	1:B:441:LEU:HD23	1.89	0.55
1:B:218:ARG:NE	1:B:218:ARG:H	2.07	0.53
1:A:450:ALA:HB1	4:A:2327:HOH:O	2.09	0.52
1:A:417:LEU:CG	1:A:441:LEU:HD21	2.40	0.51
1:B:65:HIS:O	1:B:85:GLY:HA3	2.10	0.50
1:B:466:ARG:HD3	1:B:517:TRP:CZ2	2.48	0.48
1:B:218:ARG:H	1:B:218:ARG:CD	2.26	0.48
1:A:52:SER:OG	1:A:53:PRO:HD3	2.13	0.48
1:B:52:SER:OG	1:B:53:PRO:HD3	2.16	0.45
1:A:456:THR:O	1:A:470:GLN:HA	2.16	0.45
1:A:413:THR:HG23	1:A:441:LEU:HD23	1.99	0.45
1:B:456:THR:O	1:B:470:GLN:HA	2.16	0.45
1:A:156:ILE:HD11	1:A:212:TRP:CH2	2.52	0.44
1:A:218:ARG:NE	1:A:218:ARG:H	2.14	0.44
1:A:273:ARG:NH2	4:A:2219:HOH:O	2.51	0.43
1:A:325:LEU:HD12	1:A:358:TRP:CE3	2.53	0.43
1:A:218:ARG:H	1:A:218:ARG:CD	2.32	0.43
1:B:43:THR:OG1	1:B:402:HIS:HD2	2.02	0.42
1:B:334:ASP:OD2	1:B:341[B]:ARG:NH2	2.49	0.41
1:B:484:PRO:O	1:B:485:ALA:HB3	2.22	0.40
1:B:422:HIS:CG	1:B:423:PRO:HD2	2.56	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	526/536 (98%)	512 (97%)	14 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	528/536 (98%)	514 (97%)	14 (3%)	0	100	100
All	All	1054/1072 (98%)	1026 (97%)	28 (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	402/429 (94%)	394 (98%)	8 (2%)	63	20
1	B	404/429 (94%)	396 (98%)	8 (2%)	63	20
All	All	806/858 (94%)	790 (98%)	16 (2%)	63	20

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	218	ARG
1	A	270	ASP
1	A	318	ASN
1	A	359	TYR
1	A	394	LEU
1	A	402	HIS
1	A	441	LEU
1	A	506	GLN
1	B	218	ARG
1	B	270	ASP
1	B	318	ASN
1	B	359	TYR
1	B	394	LEU
1	B	441	LEU
1	B	506	GLN
1	B	519	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	HIS
1	B	89	HIS
1	B	402	HIS
1	B	506	GLN

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	$\# Z > 2$
1	DDZ	A	51	1,2	2,6,7	0.61	0	1,7,9	3.04	1 (100%)
1	DDZ	B	51	1,2	2,6,7	0.32	0	1,7,9	3.25	1 (100%)

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	DDZ	A	51	1,2	-	0/0/6/8	0/0/0/0
1	DDZ	B	51	1,2	-	0/0/6/8	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	51	DDZ	C-CA-N	3.04	115.73	109.12
1	B	51	DDZ	C-CA-N	3.25	116.18	109.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	A	1529	2	4,4,4	0.80	0	6,6,6	0.14	0
3	SO4	A	1530	-	4,4,4	0.59	0	6,6,6	0.07	0
3	SO4	A	1531	-	4,4,4	0.60	0	6,6,6	0.14	0
3	SO4	A	1532	-	4,4,4	0.60	0	6,6,6	0.07	0
3	SO4	B	1529	2	4,4,4	0.84	0	6,6,6	0.29	0
3	SO4	B	1530	-	4,4,4	0.61	0	6,6,6	0.06	0
3	SO4	B	1531	-	4,4,4	0.64	0	6,6,6	0.23	0
3	SO4	B	1532	-	4,4,4	0.60	0	6,6,6	0.07	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
3	SO4	A	1529	2	-	0/0/0/0	0/0/0/0
3	SO4	A	1530	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1531	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1532	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1529	2	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	B	1530	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1531	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1532	-	-	0/0/0/0	0/0/0/0

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 [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	524/536 (97%)	0.43	29 (5%)	29 25	7, 13, 28, 58	3 (0%)
1	B	524/536 (97%)	0.49	30 (5%)	27 23	7, 13, 28, 58	4 (0%)
All	All	1048/1072 (97%)	0.46	59 (5%)	28 24	7, 13, 28, 58	7 (0%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	173	THR	6.0
1	B	78	PRO	5.6
1	B	77	THR	5.2
1	A	173	THR	5.0
1	A	78	PRO	4.3
1	B	172	ASP	4.0
1	B	261	LEU	3.9
1	B	482	VAL	3.7
1	A	177	GLY	3.6
1	B	481	PRO	3.6
1	A	77	THR	3.5
1	A	172	ASP	3.4
1	B	429	GLY	3.4
1	B	79	GLU	3.4
1	A	481	PRO	3.3
1	B	80	LEU	3.2
1	A	483	GLY	3.2
1	A	480	ALA	3.2
1	A	175	PRO	3.2
1	B	329	PRO	3.1
1	A	79	GLU	3.0
1	B	175	PRO	3.0
1	A	76	LEU	2.9
1	B	476	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	82	GLY	2.9
1	B	480	ALA	2.8
1	B	174	LEU	2.7
1	B	76	LEU	2.7
1	A	329	PRO	2.6
1	A	3	LYS	2.6
1	B	3	LYS	2.6
1	A	174	LEU	2.5
1	A	479	PRO	2.5
1	B	394	LEU	2.5
1	A	82	GLY	2.5
1	A	149	TYR	2.4
1	A	463	PHE	2.4
1	B	504	ASP	2.4
1	A	527	VAL	2.3
1	A	484	PRO	2.3
1	A	486	THR	2.3
1	B	441	LEU	2.3
1	A	170	TYR	2.2
1	A	169	ARG	2.2
1	A	394	LEU	2.2
1	B	188	LEU	2.2
1	A	188	LEU	2.2
1	B	251	ALA	2.2
1	B	463	PHE	2.1
1	B	479	PRO	2.1
1	B	341[A]	ARG	2.1
1	B	507	PRO	2.0
1	B	176	GLU	2.0
1	A	150	ASP	2.0
1	B	252	ASP	2.0
1	A	331	PHE	2.0
1	A	176	GLU	2.0
1	B	170	TYR	2.0
1	A	251	ALA	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(\AA^2)	Q<0.9
1	DDZ	A	51	7/8	0.95	0.09	-	9,11,12,16	0
1	DDZ	B	51	7/8	0.95	0.08	-	10,11,12,16	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	B	1529	5/5	0.90	0.11	3.13	13,17,18,21	0
3	SO4	B	1532	5/5	0.93	0.18	2.82	20,25,38,41	5
3	SO4	A	1529	5/5	0.91	0.12	2.66	12,15,17,19	0
3	SO4	B	1531	5/5	0.97	0.08	-0.52	17,18,19,28	0
3	SO4	A	1531	5/5	0.98	0.07	-1.02	18,19,21,23	0
2	CA	A	1528	1/1	0.99	0.05	-2.47	10,10,10,10	0
2	CA	B	1528	1/1	0.99	0.04	-2.89	10,10,10,10	0
3	SO4	A	1532	5/5	0.58	0.40	-	31,33,39,53	5
3	SO4	B	1530	5/5	0.94	0.20	-	27,29,35,42	0
3	SO4	A	1530	5/5	0.95	0.20	-	24,28,36,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.