



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2026 – 11:46 AM UTC

PDB ID : 1B4K / pdb_00001b4k
Title : High resolution crystal structure of a MG2-dependent 5-aminolevulinic acid dehydratase
Authors : Frankenberg, N.; Jahn, D.; Heinz, D.W.
Deposited on : 1998-12-22
Resolution : 1.67 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

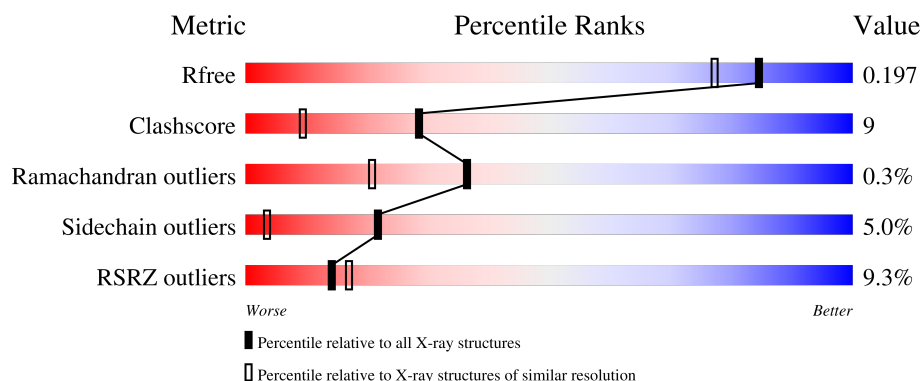
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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	1054 (1.68-1.68)
Clashscore	190562	1078 (1.68-1.68)
Ramachandran outliers	187476	1068 (1.68-1.68)
Sidechain outliers	187428	1067 (1.68-1.68)
RSRZ outliers	180081	1055 (1.68-1.68)

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	337	4% 82% 13% ••
1	B	337	14% 75% 16% •• 5%

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
4	SHF	A	403	-	X	-	-
4	SHF	B	402	-	X	-	-

2 Entry composition [i](#)

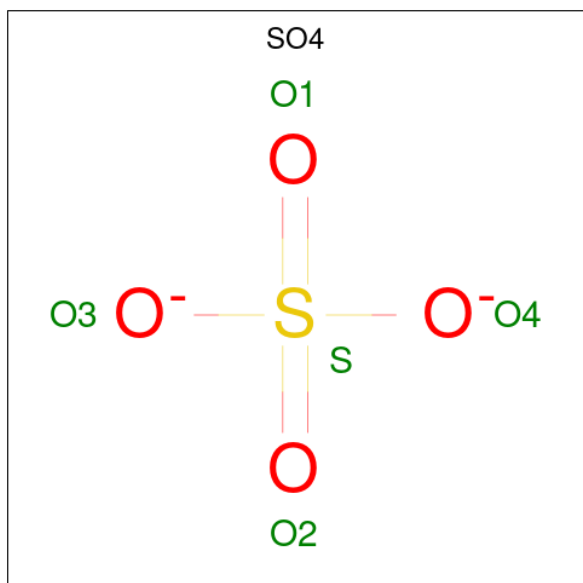
There are 5 unique types of molecules in this entry. The entry contains 5604 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (5-AMINOLEVULINIC ACID DEHYDRATASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	0	0
			2522	1586	443	483	10			
1	B	320	Total	C	N	O	S	0	0	0
			2485	1563	437	475	10			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

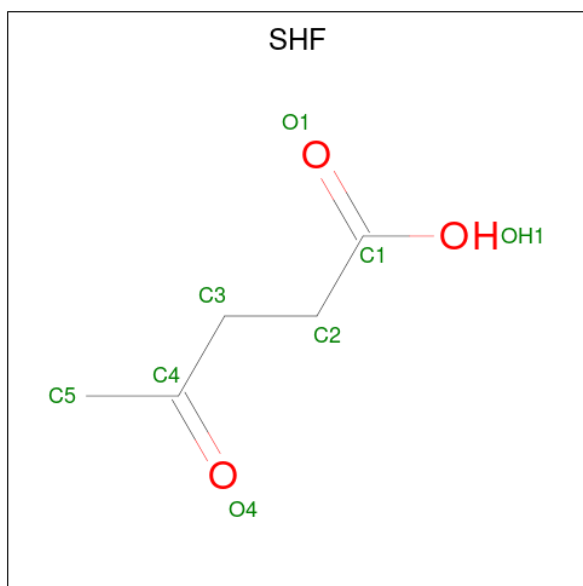


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is LAEVULINIC ACID (CCD ID: SHF) (formula: $C_5H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	5	2		
4	B	1	Total	C	O	0	0
			7	5	2		

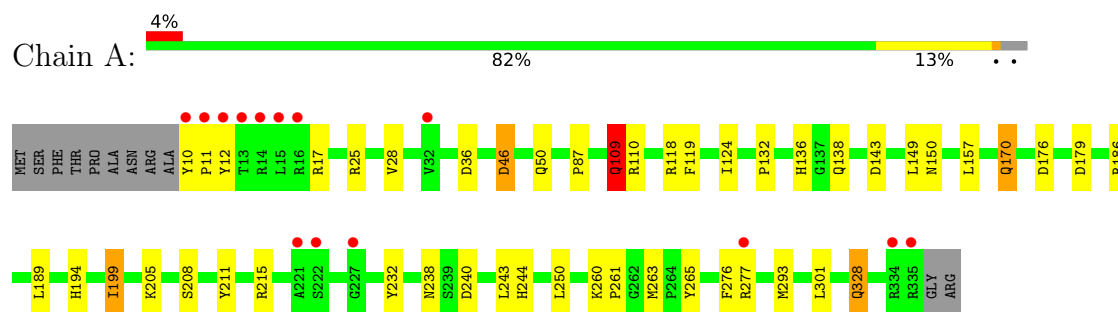
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	322	Total	O	0	0
			322	322		
5	B	250	Total	O	0	0
			250	250		

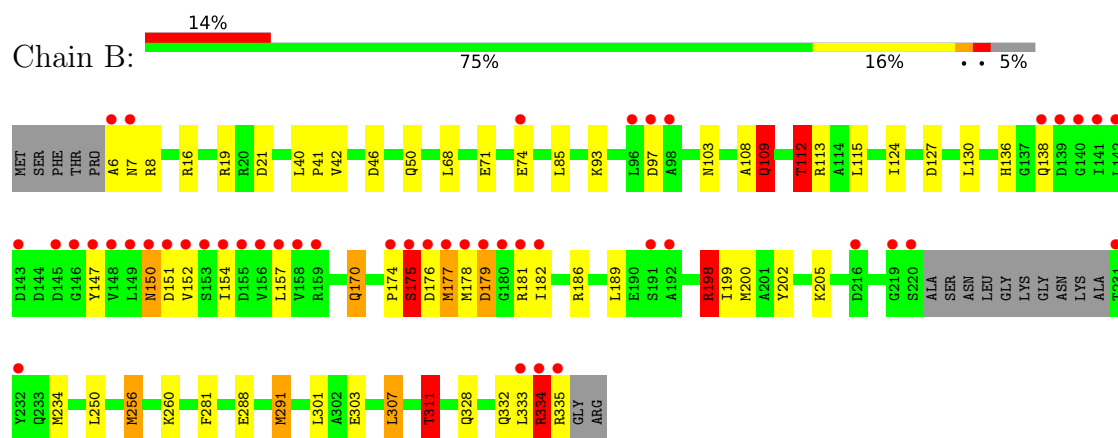
3 Residue-property plots [i](#)

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: PROTEIN (5-AMINOLEVULINIC ACID DEHYDRATASE)



• Molecule 1: PROTEIN (5-AMINOLEVULINIC ACID DEHYDRATASE)



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	128.24Å 128.24Å 86.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.30 – 1.67 51.30 – 1.67	Depositor EDS
% Data completeness (in resolution range)	98.2 (51.30-1.67) 95.3 (51.30-1.67)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 1.67Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.178 , 0.208 0.170 , 0.197	Depositor DCC
R_{free} test set	4105 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5604	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.15% 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: SHF, SO4, MG

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	0.71	0/2569	1.44	22/3486 (0.6%)
1	B	0.67	0/2531	1.52	25/3435 (0.7%)
All	All	0.69	0/5100	1.48	47/6921 (0.7%)

There are no bond length outliers.

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	113	ARG	CD-NE-CZ	23.39	157.14	124.40
1	B	198	ARG	CD-NE-CZ	18.30	150.01	124.40
1	A	25	ARG	CD-NE-CZ	10.66	139.32	124.40
1	A	118	ARG	CD-NE-CZ	9.10	137.14	124.40
1	B	150	ASN	N-CA-C	8.57	123.03	112.23
1	A	215	ARG	CD-NE-CZ	8.21	135.90	124.40
1	B	179	ASP	CA-CB-CG	7.91	120.51	112.60
1	B	152	VAL	N-CA-C	-7.80	103.19	110.53
1	A	25	ARG	NE-CZ-NH2	-7.62	112.34	119.20
1	B	175	SER	CA-C-O	7.54	128.40	119.28
1	B	19	ARG	NE-CZ-NH2	-7.40	112.54	119.20
1	B	198	ARG	NE-CZ-NH1	-7.17	114.33	121.50
1	A	328	GLN	OE1-CD-NE2	6.83	129.44	122.60
1	A	10	TYR	CA-C-O	-6.40	109.92	120.80
1	B	291	MET	CB-CA-C	-6.12	101.28	110.88
1	A	46	ASP	CA-CB-CG	6.06	118.66	112.60
1	B	97	ASP	CA-CB-CG	5.92	118.52	112.60
1	A	194	HIS	CA-C-N	5.92	128.21	120.28
1	A	194	HIS	C-N-CA	5.92	128.21	120.28
1	B	281	PHE	CA-CB-CG	-5.82	107.98	113.80
1	A	10	TYR	CB-CG-CD1	-5.77	112.14	120.80
1	B	147	TYR	CA-CB-CG	-5.75	103.55	113.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	303	GLU	CB-CG-CD	-5.71	102.90	112.60
1	A	10	TYR	CA-CB-CG	-5.64	103.74	113.90
1	B	174	PRO	CB-CA-C	5.63	120.86	111.56
1	A	176	ASP	CA-CB-CG	-5.62	106.98	112.60
1	A	36	ASP	CA-CB-CG	5.61	118.21	112.60
1	A	232	TYR	CA-CB-CG	5.47	123.75	113.90
1	B	46	ASP	CA-C-N	-5.40	115.83	122.75
1	B	46	ASP	C-N-CA	-5.40	115.83	122.75
1	B	7	ASN	CA-CB-CG	5.37	117.97	112.60
1	A	176	ASP	N-CA-C	5.34	118.96	112.23
1	A	238	ASN	N-CA-C	5.33	117.93	109.24
1	A	263	MET	CG-SD-CE	-5.32	89.21	100.90
1	B	103	ASN	CA-C-N	5.31	125.03	119.82
1	B	103	ASN	C-N-CA	5.31	125.03	119.82
1	A	10	TYR	CB-CG-CD2	5.31	128.77	120.80
1	B	291	MET	CG-SD-CE	5.29	112.53	100.90
1	A	109	GLN	OE1-CD-NE2	-5.28	117.32	122.60
1	B	198	ARG	NE-CZ-NH2	5.26	123.93	119.20
1	B	19	ARG	NE-CZ-NH1	5.21	126.72	121.50
1	B	109	GLN	OE1-CD-NE2	-5.17	117.43	122.60
1	B	112	THR	N-CA-CB	5.14	117.77	110.16
1	A	119	PHE	CA-C-N	5.13	125.42	119.47
1	A	119	PHE	C-N-CA	5.13	125.42	119.47
1	A	208	SER	N-CA-C	5.09	117.90	109.85
1	B	311	THR	N-CA-CB	5.09	118.18	110.28

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	2522	0	2490	38	0
1	B	2485	0	2449	51	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
4	A	7	0	7	0	0
4	B	7	0	7	1	0
5	A	322	0	0	8	0
5	B	250	0	0	5	0
All	All	5604	0	4953	88	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 9.

All (88) 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)
4:B:402:SHF:H51	5:B:690:HOH:O	1.48	1.11
1:A:150:ASN:HD21	1:A:179:ASP:H	1.19	0.87
1:B:175:SER:HA	1:B:202:TYR:CD1	2.14	0.82
1:A:170:GLN:H	1:A:170:GLN:HE21	1.28	0.81
1:A:189:LEU:CD1	1:A:199:ILE:HD11	2.10	0.80
1:A:136:HIS:HD2	1:A:138:GLN:H	1.30	0.79
1:A:189:LEU:HD12	1:A:199:ILE:HD11	1.64	0.79
1:B:170:GLN:HE21	1:B:170:GLN:H	1.36	0.74
1:A:186:ARG:NH1	1:A:199:ILE:HD13	2.04	0.73
1:B:136:HIS:HD2	1:B:138:GLN:H	1.35	0.73
1:B:198:ARG:N	1:B:198:ARG:HD3	2.02	0.73
1:B:200:MET:HA	1:B:256:MET:HG2	1.74	0.70
1:A:186:ARG:CZ	1:A:199:ILE:HD13	2.23	0.69
1:A:150:ASN:HD21	1:A:179:ASP:N	1.90	0.68
1:A:87:PRO:HD2	5:A:637:HOH:O	1.93	0.68
1:B:176:ASP:OD2	1:B:181:ARG:HD2	1.93	0.68
1:B:205:LYS:HE3	1:B:260:LYS:HE2	1.76	0.67
1:A:205:LYS:HE3	1:A:260:LYS:HE2	1.77	0.66
1:B:186:ARG:CZ	1:B:199:ILE:HD13	2.25	0.66
1:B:307:LEU:O	1:B:311:THR:HG23	1.95	0.66
1:B:189:LEU:HD12	1:B:199:ILE:HD11	1.79	0.64
1:B:108:ALA:O	1:B:112:THR:HG23	1.99	0.63
1:B:157:LEU:HD13	1:B:181:ARG:HD3	1.81	0.62
1:B:109:GLN:HE21	1:B:109:GLN:H	1.48	0.61
1:A:150:ASN:HD22	1:B:16:ARG:HH11	1.47	0.61
1:B:124:ILE:H	1:B:170:GLN:NE2	1.99	0.61
1:B:178:MET:HG3	1:B:181:ARG:CZ	2.31	0.61
1:B:189:LEU:CD1	1:B:199:ILE:HD11	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:ARG:NH2	1:A:199:ILE:HD13	2.18	0.59
1:B:177:MET:HE2	1:B:205:LYS:H	1.68	0.59
1:A:136:HIS:CD2	1:A:138:GLN:H	2.17	0.59
1:B:40:LEU:HD23	1:B:41:PRO:O	2.03	0.59
1:B:186:ARG:NH2	1:B:199:ILE:HD13	2.19	0.58
1:A:328:GLN:HG3	5:A:735:HOH:O	2.04	0.57
1:A:132:PRO:HG3	5:A:784:HOH:O	2.04	0.57
1:B:136:HIS:CD2	1:B:138:GLN:H	2.20	0.56
1:B:109:GLN:H	1:B:109:GLN:NE2	2.03	0.55
1:B:182:ILE:HD13	5:B:723:HOH:O	2.06	0.55
1:A:124:ILE:H	1:A:170:GLN:NE2	2.05	0.54
1:A:109:GLN:HE21	1:A:109:GLN:H	1.56	0.54
1:A:243:LEU:HD22	1:A:243:LEU:H	1.73	0.54
1:B:200:MET:HG3	1:B:256:MET:HG3	1.90	0.53
1:B:112:THR:HB	1:B:124:ILE:HG21	1.90	0.53
1:B:124:ILE:H	1:B:170:GLN:HE22	1.55	0.52
1:A:170:GLN:HE21	1:A:170:GLN:N	2.04	0.52
1:B:40:LEU:HD22	1:B:42:VAL:CG1	2.39	0.52
1:A:136:HIS:HE1	5:A:527:HOH:O	1.94	0.51
1:B:6:ALA:HA	5:B:696:HOH:O	2.11	0.50
1:A:109:GLN:H	1:A:109:GLN:NE2	2.09	0.50
1:A:186:ARG:HH12	1:A:199:ILE:HD13	1.75	0.50
1:B:334:ARG:O	1:B:335:ARG:C	2.54	0.50
1:B:177:MET:HG2	1:B:205:LYS:HB3	1.94	0.49
1:A:240:ASP:HA	1:A:243:LEU:HD23	1.94	0.49
1:B:178:MET:HE2	1:B:181:ARG:HH21	1.77	0.48
1:B:288:GLU:HA	1:B:291:MET:HE2	1.94	0.48
1:A:276:PHE:O	1:A:277:ARG:HG2	2.13	0.47
1:B:130:LEU:HG	1:B:157:LEU:HD21	1.96	0.47
1:B:332:GLN:O	1:B:334:ARG:N	2.47	0.47
1:A:205:LYS:CE	1:A:260:LYS:HE2	2.42	0.47
1:A:243:LEU:HD22	1:A:243:LEU:N	2.30	0.47
1:B:136:HIS:CD2	1:B:138:GLN:HB2	2.50	0.47
1:B:85:LEU:HD12	1:B:112:THR:HG22	1.96	0.46
1:B:178:MET:HG3	1:B:181:ARG:NE	2.30	0.46
1:B:127:ASP:OD1	1:B:175:SER:HB3	2.15	0.46
1:B:68:LEU:HD21	1:B:115:LEU:HD11	1.98	0.45
1:A:124:ILE:H	1:A:170:GLN:HE22	1.62	0.45
1:B:40:LEU:HD23	1:B:40:LEU:C	2.42	0.45
1:B:21:ASP:HA	5:B:512:HOH:O	2.17	0.45
1:A:261:PRO:HD2	1:A:265:TYR:CE2	2.53	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:GLN:HE21	1:B:170:GLN:N	2.11	0.44
1:B:178:MET:HG3	1:B:181:ARG:NH2	2.31	0.44
1:B:234:MET:HE2	1:B:234:MET:HB3	1.87	0.44
1:B:328:GLN:HG2	5:B:749:HOH:O	2.18	0.44
1:B:175:SER:HA	1:B:202:TYR:CG	2.52	0.44
1:A:143:ASP:HB3	1:A:149:LEU:HD21	2.00	0.43
1:A:211:TYR:CE1	1:A:260:LYS:HE3	2.53	0.43
1:A:244:HIS:CD2	1:B:8:ARG:HH11	2.37	0.43
1:A:28:VAL:HG13	5:A:623:HOH:O	2.19	0.42
1:B:71:GLU:HA	1:B:74:GLU:OE1	2.19	0.42
1:A:46:ASP:OD1	1:A:110:ARG:NH2	2.48	0.42
1:A:211:TYR:HB2	5:A:590:HOH:O	2.19	0.42
1:B:186:ARG:NH1	1:B:199:ILE:HD13	2.35	0.42
1:B:40:LEU:HD22	1:B:42:VAL:HG13	2.02	0.42
1:A:12:TYR:HB2	5:A:749:HOH:O	2.19	0.42
1:A:150:ASN:ND2	1:A:179:ASP:H	2.01	0.41
1:A:11:PRO:O	1:A:17:ARG:NH2	2.54	0.41
1:A:277:ARG:HB3	5:A:736:HOH:O	2.20	0.40
1:B:175:SER:HB2	1:B:202:TYR:CE1	2.56	0.40

There are no symmetry-related clashes.

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	324/337 (96%)	317 (98%)	7 (2%)	0	100	100
1	B	316/337 (94%)	307 (97%)	7 (2%)	2 (1%)	21	7
All	All	640/674 (95%)	624 (98%)	14 (2%)	2 (0%)	36	21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	333	LEU
1	B	334	ARG

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	261/269 (97%)	253 (97%)	8 (3%)	35	10
1	B	257/269 (96%)	239 (93%)	18 (7%)	14	1
All	All	518/538 (96%)	492 (95%)	26 (5%)	22	3

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	109	GLN
1	A	157	LEU
1	A	170	GLN
1	A	199	ILE
1	A	250	LEU
1	A	293	MET
1	A	301	LEU
1	B	50	GLN
1	B	93	LYS
1	B	109	GLN
1	B	112	THR
1	B	150	ASN
1	B	151	ASP
1	B	154	ILE
1	B	170	GLN
1	B	175	SER
1	B	177	MET
1	B	179	ASP
1	B	198	ARG
1	B	250	LEU
1	B	256	MET

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Mol	Chain	Res	Type
1	B	301	LEU
1	B	307	LEU
1	B	311	THR
1	B	334	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	109	GLN
1	A	136	HIS
1	A	150	ASN
1	A	170	GLN
1	A	244	HIS
1	B	109	GLN
1	B	136	HIS
1	B	170	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](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	$\# Z > 2$
2	SO4	B	401	-	4,4,4	0.73	0	6,6,6	0.11	0
4	SHF	B	402	1	6,6,7	1.80	1 (16%)	6,6,8	3.87	4 (66%)
4	SHF	A	403	1	6,6,7	2.07	1 (16%)	6,6,8	3.39	4 (66%)
2	SO4	A	401	-	4,4,4	0.68	0	6,6,6	0.38	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	SHF	B	402	1	-	2/4/4/5	-
4	SHF	A	403	1	-	2/4/4/5	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	403	SHF	O1-C1	4.05	1.35	1.22
4	B	402	SHF	O1-C1	3.46	1.33	1.22

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	402	SHF	C3-C2-C1	5.90	129.90	114.51
4	B	402	SHF	OH1-C1-C2	5.65	131.85	114.00
4	A	403	SHF	C3-C2-C1	5.21	128.10	114.51
4	A	403	SHF	OH1-C1-C2	4.77	129.06	114.00
4	B	402	SHF	O1-C1-C2	-4.04	110.28	123.09
4	A	403	SHF	O1-C1-C2	-3.28	112.69	123.09
4	A	403	SHF	OH1-C1-O1	-2.68	116.43	123.33
4	B	402	SHF	OH1-C1-O1	-2.37	117.24	123.33

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	403	SHF	C2-C3-C4-C5

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Mol	Chain	Res	Type	Atoms
4	B	402	SHF	C2-C3-C4-C5
4	B	402	SHF	OH1-C1-C2-C3
4	A	403	SHF	OH1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	402	SHF	1	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	326/337 (96%)	-0.12	14 (4%) 40 48	11, 16, 30, 76	0
1	B	320/337 (94%)	0.42	46 (14%) 6 7	12, 20, 54, 68	0
All	All	646/674 (95%)	0.14	60 (9%) 14 17	11, 18, 44, 76	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	TYR	7.2
1	B	6	ALA	6.5
1	A	10	TYR	6.1
1	B	149	LEU	6.0
1	B	175	SER	5.7
1	B	147	TYR	5.5
1	A	11	PRO	5.1
1	B	148	VAL	5.0
1	B	176	ASP	5.0
1	B	177	MET	5.0
1	A	277	ARG	4.9
1	B	335	ARG	4.6
1	A	13	THR	4.5
1	B	334	ARG	4.4
1	B	152	VAL	4.4
1	B	180	GLY	4.2
1	B	179	ASP	4.1
1	B	142	LEU	4.1
1	B	141	ILE	4.1
1	A	15	LEU	3.9
1	B	156	VAL	3.7
1	B	150	ASN	3.6
1	B	151	ASP	3.5
1	A	335	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	181	ARG	3.3
1	B	174	PRO	3.3
1	B	220	SER	3.2
1	B	182	ILE	3.2
1	B	178	MET	3.1
1	A	221	ALA	3.1
1	B	143	ASP	3.1
1	A	222	SER	3.1
1	B	219	GLY	3.0
1	B	96	LEU	3.0
1	B	157	LEU	3.0
1	B	154	ILE	2.9
1	A	14	ARG	2.9
1	B	98	ALA	2.9
1	B	74	GLU	2.8
1	A	16	ARG	2.8
1	B	145	ASP	2.8
1	B	155	ASP	2.8
1	B	232	TYR	2.8
1	B	146	GLY	2.8
1	B	153	SER	2.7
1	A	227	GLY	2.7
1	B	333	LEU	2.7
1	B	140	GLY	2.6
1	B	231	THR	2.6
1	A	334	ARG	2.4
1	B	192	ALA	2.4
1	B	158	VAL	2.3
1	B	7	ASN	2.3
1	B	97	ASP	2.2
1	B	191	SER	2.2
1	A	32	VAL	2.1
1	B	138	GLN	2.1
1	B	139	ASP	2.1
1	B	159	ARG	2.1
1	B	216	ASP	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

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($\text{\AA}^2$)	Q<0.9
2	SO4	A	401	5/5	0.74	0.20	52,53,54,54	0
2	SO4	B	401	5/5	0.81	0.17	76,76,76,76	0
4	SHF	A	403	7/8	0.87	0.11	14,15,17,17	0
4	SHF	B	402	7/8	0.88	0.12	17,17,23,24	0
3	MG	A	402	1/1	0.99	0.02	12,12,12,12	0

6.5 Other polymers [i](#)

There are no such residues in this entry.