



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:40 PM UTC

PDB ID : 6HSR / pdb_00006hsr
Title : The crystal structure of type II Dehydroquinase from *Psychromonas ingrahamii* 37, 40% ethanol as cryoprotectant
Authors : Lapthorn, A.J.; Roszak, A.W.
Deposited on : 2018-10-01
Resolution : 2.00 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:

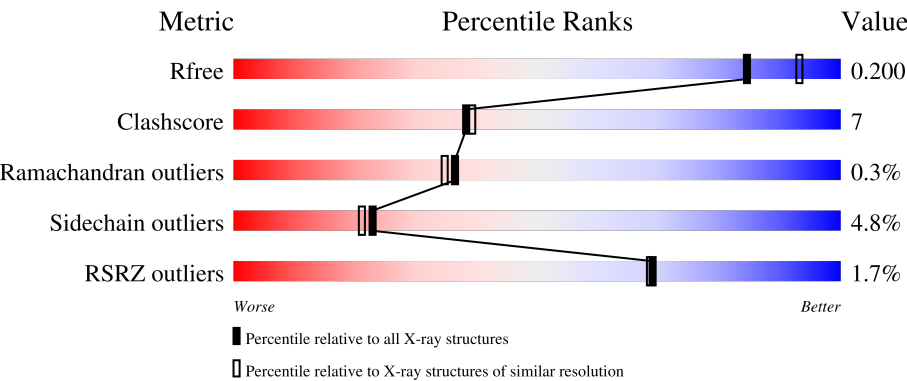
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (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 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	152	2% 76% 18% ...
1	B	152	2% 66% 30% ...
1	C	152	2% 79% 16% ...
1	D	152	% 68% 24% 7% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4906 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 3-dehydroquinate dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	148	Total	C	N	O	S	0	1	0
			1164	742	205	216	1			
1	B	151	Total	C	N	O	S	0	1	0
			1179	751	208	219	1			
1	C	149	Total	C	N	O	S	0	0	0
			1158	737	203	217	1			
1	D	152	Total	C	N	O	S	0	1	0
			1189	756	209	222	2			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP A1SZA3
A	-1	SER	-	expression tag	UNP A1SZA3
A	0	HIS	-	expression tag	UNP A1SZA3
B	-2	GLY	-	expression tag	UNP A1SZA3
B	-1	SER	-	expression tag	UNP A1SZA3
B	0	HIS	-	expression tag	UNP A1SZA3
C	-2	GLY	-	expression tag	UNP A1SZA3
C	-1	SER	-	expression tag	UNP A1SZA3
C	0	HIS	-	expression tag	UNP A1SZA3
D	-2	GLY	-	expression tag	UNP A1SZA3
D	-1	SER	-	expression tag	UNP A1SZA3
D	0	HIS	-	expression tag	UNP A1SZA3

- Molecule 2 is ETHANOL (CCD ID: EOH) (formula: C₂H₆O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			3	2	1		
2	B	1	Total	C	O	0	0
			3	2	1		
2	C	1	Total	C	O	0	0
			3	2	1		
2	D	1	Total	C	O	0	0
			3	2	1		

- Molecule 3 is SULFATE ION (CCD ID: 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	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	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

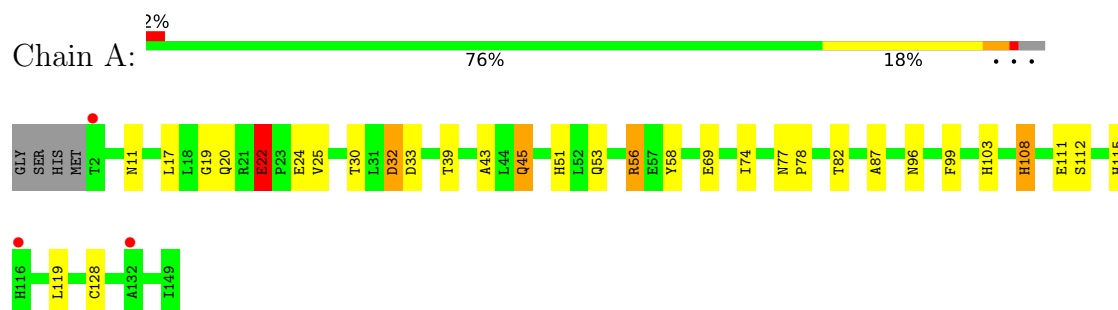
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	20	Total O 20 20	0	0
4	B	42	Total O 42 42	0	0
4	C	16	Total O 16 16	0	0
4	D	76	Total O 76 76	0	0

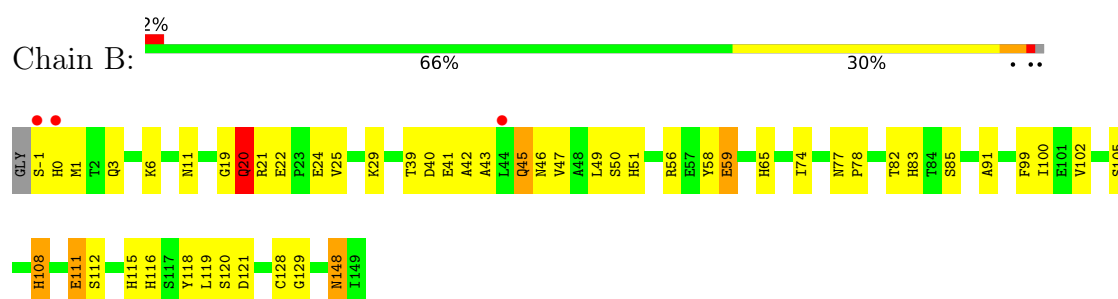
3 Residue-property plots [i](#)

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.

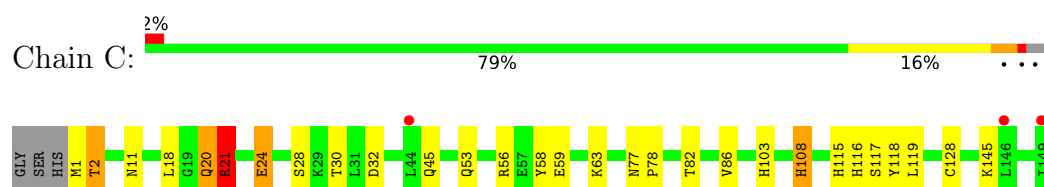
- Molecule 1: 3-dehydroquinase dehydratase



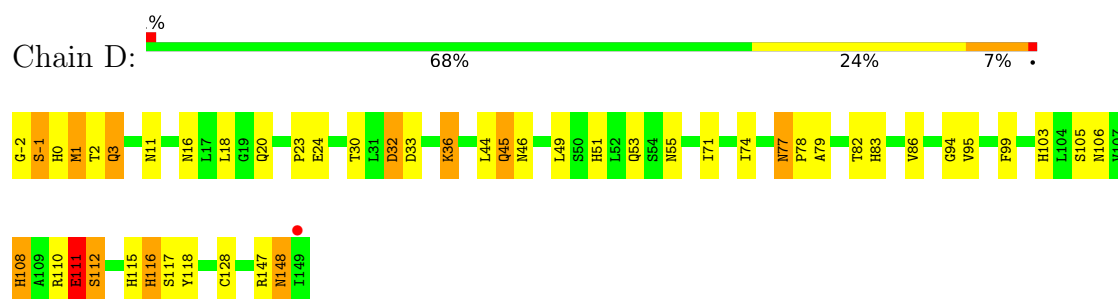
- Molecule 1: 3-dehydroquinase dehydratase



- Molecule 1: 3-dehydroquinase dehydratase



- Molecule 1: 3-dehydroquinase dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	138.83Å 138.83Å 138.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	69.51 – 2.00 69.51 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.5 (69.51-2.00) 94.5 (69.51-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.173 , 0.194 0.181 , 0.200	Depositor DCC
R_{free} test set	2836 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	54.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.036 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4906	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	1.53	14/1188 (1.2%)	1.58	7/1608 (0.4%)
1	B	1.83	23/1203 (1.9%)	1.77	17/1629 (1.0%)
1	C	1.48	7/1178 (0.6%)	1.56	9/1596 (0.6%)
1	D	2.00	23/1213 (1.9%)	1.84	22/1641 (1.3%)
All	All	1.72	67/4782 (1.4%)	1.69	55/6474 (0.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	D	0	4
All	All	0	6

All (67) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	108	HIS	CE1-NE2	18.03	1.50	1.32
1	B	108	HIS	CE1-NE2	15.95	1.48	1.32
1	B	41	GLU	C-O	-11.09	1.11	1.24
1	D	111	GLU	CD-OE2	-10.60	1.05	1.25
1	D	95	VAL	C-O	-10.26	1.12	1.24
1	C	108	HIS	CE1-NE2	9.70	1.42	1.32
1	A	103	HIS	CE1-NE2	9.39	1.42	1.32
1	D	117	SER	CA-CB	-9.18	1.32	1.53
1	A	115	HIS	C-O	-9.06	1.12	1.24
1	D	103	HIS	CG-ND1	9.04	1.48	1.38
1	D	112	SER	CA-CB	8.95	1.69	1.53

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	51	HIS	CE1-NE2	8.63	1.41	1.32
1	D	115	HIS	C-O	-8.28	1.13	1.24
1	D	116	HIS	CG-ND1	7.83	1.46	1.38
1	A	87	ALA	C-O	7.80	1.33	1.24
1	A	108	HIS	CE1-NE2	7.72	1.40	1.32
1	A	51	HIS	C-O	7.71	1.32	1.23
1	A	43	ALA	C-O	7.67	1.33	1.24
1	A	56	ARG	C-O	7.58	1.32	1.24
1	B	42	ALA	N-CA	7.56	1.55	1.46
1	C	116	HIS	CG-ND1	7.44	1.46	1.38
1	B	115	HIS	CG-ND1	7.43	1.46	1.38
1	D	94	GLY	C-O	-7.30	1.15	1.23
1	B	39	THR	CA-C	7.21	1.62	1.52
1	A	111	GLU	C-O	7.20	1.32	1.23
1	B	83	HIS	C-O	-7.14	1.14	1.24
1	C	116	HIS	CE1-NE2	6.84	1.39	1.32
1	D	49	LEU	C-O	-6.80	1.15	1.24
1	B	115	HIS	C-O	-6.75	1.15	1.24
1	D	105	SER	CB-OG	6.69	1.55	1.42
1	C	115	HIS	C-O	-6.68	1.15	1.24
1	D	71	ILE	C-O	-6.65	1.17	1.24
1	D	86	VAL	C-O	6.46	1.32	1.24
1	C	86	VAL	C-O	6.46	1.32	1.24
1	C	103	HIS	CE1-NE2	6.40	1.39	1.32
1	C	117	SER	CA-CB	-6.38	1.42	1.53
1	A	103	HIS	CG-ND1	6.37	1.45	1.38
1	B	43	ALA	C-O	6.30	1.31	1.24
1	B	50	SER	C-O	6.29	1.32	1.23
1	D	46	ASN	C-O	6.18	1.31	1.23
1	B	108	HIS	C-O	6.15	1.31	1.24
1	B	129	GLY	C-O	-6.12	1.15	1.24
1	B	91	ALA	C-O	6.08	1.31	1.24
1	D	116	HIS	N-CA	6.05	1.53	1.46
1	D	36	LYS	C-O	6.04	1.31	1.24
1	D	32	ASP	CG-OD2	5.85	1.36	1.25
1	B	105	SER	N-CA	5.76	1.52	1.45
1	B	108	HIS	CD2-NE2	5.75	1.44	1.37
1	B	46	ASN	C-N	5.73	1.40	1.33
1	D	33	ASP	C-O	-5.67	1.17	1.24
1	B	65	HIS	ND1-CE1	-5.62	1.26	1.32
1	B	85	SER	CA-CB	-5.62	1.44	1.52
1	D	106	ASN	C-O	5.60	1.30	1.24

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	102	VAL	C-O	5.50	1.29	1.24
1	A	33	ASP	C-O	-5.44	1.17	1.24
1	D	18	LEU	C-O	5.39	1.30	1.23
1	A	39	THR	C-O	5.23	1.30	1.24
1	B	6	LYS	CG-CD	-5.23	1.36	1.52
1	B	20	GLN	C-O	-5.22	1.17	1.23
1	B	51	HIS	CE1-NE2	5.20	1.37	1.32
1	B	100	ILE	C-O	-5.18	1.18	1.24
1	A	111	GLU	CD-OE2	5.17	1.35	1.25
1	A	69	GLU	C-O	-5.05	1.15	1.24
1	A	51	HIS	CE1-NE2	5.04	1.37	1.32
1	D	83	HIS	CE1-NE2	5.03	1.37	1.32
1	D	115	HIS	CA-C	5.02	1.59	1.52
1	B	41	GLU	CA-C	5.00	1.59	1.52

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	45	GLN	CB-CG-CD	-13.82	89.11	112.60
1	D	108	HIS	CE1-NE2-CD2	-9.32	99.68	109.00
1	B	108	HIS	CE1-NE2-CD2	-8.79	100.20	109.00
1	A	45	GLN	CB-CG-CD	-8.47	98.20	112.60
1	D	103	HIS	ND1-CE1-NE2	8.28	116.68	108.40
1	B	40	ASP	CA-C-O	-7.82	112.61	120.82
1	B	41	GLU	N-CA-C	-7.78	102.75	111.07
1	B	46	ASN	CA-CB-CG	-7.72	104.88	112.60
1	D	23	PRO	CA-C-N	7.57	131.17	120.28
1	D	23	PRO	C-N-CA	7.57	131.17	120.28
1	A	53	GLN	CA-C-O	-7.33	112.94	120.71
1	B	0	HIS	CA-C-N	7.24	133.65	122.42
1	B	0	HIS	C-N-CA	7.24	133.65	122.42
1	B	120	SER	CA-CB-OG	-7.20	96.71	111.10
1	B	20	GLN	CB-CA-C	-7.19	97.40	110.56
1	D	3	GLN	N-CA-CB	7.01	120.39	110.36
1	A	32	ASP	CA-CB-CG	-6.55	106.05	112.60
1	B	41	GLU	CB-CA-C	6.44	120.98	110.88
1	D	116	HIS	CA-CB-CG	-6.32	107.48	113.80
1	D	112	SER	CA-CB-OG	-6.28	98.54	111.10
1	C	116	HIS	CB-CA-C	6.27	120.09	109.75
1	D	32	ASP	CA-CB-CG	-6.21	106.39	112.60
1	A	30	THR	CA-CB-OG1	-6.20	100.30	109.60
1	B	46	ASN	OD1-CG-ND2	-6.10	116.50	122.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	40	ASP	O-C-N	6.09	128.34	122.07
1	C	118	TYR	N-CA-C	-5.98	105.74	113.16
1	D	148	ASN	CA-C-O	-5.87	114.73	121.54
1	D	112	SER	N-CA-CB	-5.86	100.67	110.39
1	D	103	HIS	CG-ND1-CE1	-5.85	99.36	109.30
1	D	55	ASN	CA-CB-CG	-5.83	106.77	112.60
1	C	145	LYS	N-CA-C	-5.74	104.94	111.14
1	D	53	GLN	CA-C-O	-5.72	114.64	120.71
1	D	108	HIS	CG-CD2-NE2	5.66	112.86	107.20
1	C	20	GLN	CB-CG-CD	-5.61	103.06	112.60
1	D	77	ASN	OD1-CG-ND2	5.58	128.18	122.60
1	D	118	TYR	N-CA-C	-5.56	106.47	113.20
1	D	16	ASN	OD1-CG-ND2	-5.55	117.05	122.60
1	C	30	THR	CA-CB-OG1	-5.50	101.35	109.60
1	C	53	GLN	CA-C-O	-5.50	114.95	121.16
1	D	30	THR	CA-CB-OG1	-5.50	101.36	109.60
1	B	45	GLN	CB-CG-CD	-5.49	103.26	112.60
1	A	24	GLU	CB-CG-CD	5.38	121.75	112.60
1	D	83	HIS	CA-CB-CG	-5.32	108.48	113.80
1	B	148	ASN	CA-CB-CG	-5.30	107.30	112.60
1	C	108	HIS	CE1-NE2-CD2	-5.29	103.71	109.00
1	B	59	GLU	N-CA-C	-5.29	105.69	111.82
1	B	21	ARG	CG-CD-NE	-5.27	100.41	112.00
1	C	21	ARG	CB-CG-CD	5.26	123.41	111.30
1	D	115	HIS	CA-CB-CG	-5.15	108.65	113.80
1	B	108	HIS	CG-CD2-NE2	5.13	112.33	107.20
1	D	147	ARG	CG-CD-NE	-5.13	100.72	112.00
1	C	108	HIS	N-CA-C	-5.07	106.35	112.54
1	A	19	GLY	CA-C-O	-5.05	113.41	119.07
1	A	22	GLU	CB-CA-C	5.03	115.38	111.00
1	B	121	ASP	CA-C-O	-5.03	113.19	119.38

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	SER	Mainchain
1	B	111	GLU	Mainchain
1	D	110	ARG	Mainchain
1	D	111	GLU	Mainchain
1	D	112	SER	Mainchain
1	D	116	HIS	Mainchain

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	1164	0	1178	14	0
1	B	1179	0	1187	23	0
1	C	1158	0	1165	21	0
1	D	1189	0	1205	14	0
2	A	3	0	6	0	0
2	B	3	0	6	1	0
2	C	3	0	6	0	0
2	D	3	0	6	1	0
3	A	10	0	0	0	0
3	B	15	0	0	2	0
3	C	10	0	0	0	0
3	D	15	0	0	0	0
4	A	20	0	0	2	0
4	B	42	0	0	5	0
4	C	16	0	0	4	0
4	D	76	0	0	3	0
All	All	4906	0	4759	66	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 7.

All (66) 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:B:20:GLN:HE21	1:B:20:GLN:CA	1.36	1.22
1:B:20:GLN:HE21	1:B:20:GLN:N	1.52	1.06
1:B:20:GLN:HE21	1:B:20:GLN:HA	1.10	1.06
1:D:128:CYS:SG	4:D:302:HOH:O	2.14	1.04
1:B:20:GLN:CA	1:B:20:GLN:NE2	2.14	1.03
1:B:20:GLN:HA	1:B:20:GLN:NE2	1.68	0.99
1:B:128:CYS:SG	4:B:304:HOH:O	2.23	0.96
1:D:1:MET:HE3	1:D:2:THR:O	1.70	0.91
1:C:128:CYS:SG	4:C:301:HOH:O	2.36	0.83
1:A:128:CYS:SG	4:A:301:HOH:O	2.37	0.82
1:C:21:ARG:NH2	1:C:21:ARG:HG2	1.93	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:GLU:OE2	4:B:301:HOH:O	2.01	0.78
1:B:20:GLN:N	1:B:20:GLN:NE2	2.29	0.77
1:C:108:HIS:NE2	4:C:301:HOH:O	2.20	0.75
1:D:111:GLU:OE2	4:D:301:HOH:O	2.07	0.71
1:B:108:HIS:NE2	4:B:304:HOH:O	2.24	0.70
1:C:21:ARG:HG2	1:C:21:ARG:HH21	1.56	0.69
1:D:-2:GLY:O	1:D:0:HIS:N	2.26	0.68
1:D:108:HIS:NE2	4:D:302:HOH:O	2.29	0.65
2:B:201:EOH:H21	3:B:202:SO4:O3	2.02	0.60
1:C:11:ASN:HD22	1:C:77:ASN:HB3	1.68	0.59
1:C:21:ARG:HH21	1:C:21:ARG:CG	2.16	0.58
1:B:-1:SER:C	1:B:1:MET:H	2.13	0.56
1:B:29:LYS:CE	3:B:203:SO4:O1	2.56	0.54
1:C:24:GLU:H	1:C:24:GLU:CD	2.16	0.54
1:A:17:LEU:O	1:A:20:GLN:HG2	2.09	0.52
1:C:32:ASP:HB2	4:C:302:HOH:O	2.09	0.52
1:A:11:ASN:HD22	1:A:77:ASN:HB3	1.76	0.51
1:B:11:ASN:HD22	1:B:77:ASN:HB3	1.75	0.50
1:C:24:GLU:OE2	1:C:24:GLU:N	2.37	0.50
1:A:56:ARG:HG2	1:B:58:TYR:HB2	1.93	0.50
1:D:-1:SER:O	1:D:0:HIS:ND1	2.45	0.49
1:B:19:GLY:C	1:B:20:GLN:HE21	2.20	0.47
1:B:56:ARG:HG2	1:C:58:TYR:HB2	1.95	0.47
1:B:22:GLU:HB3	1:B:25:VAL:HB	1.95	0.47
1:A:108:HIS:NE2	4:A:301:HOH:O	2.36	0.47
1:C:78:PRO:HG3	1:C:119:LEU:HD12	1.96	0.46
1:A:22:GLU:HB3	1:A:25:VAL:HB	1.97	0.46
1:A:78:PRO:HG3	1:A:119:LEU:HD12	1.98	0.46
1:C:11:ASN:ND2	1:C:77:ASN:HB3	2.31	0.45
1:C:1:MET:C	1:C:2:THR:O	2.59	0.45
1:D:11:ASN:HD22	1:D:77:ASN:HB3	1.81	0.45
1:B:59:GLU:OE1	4:B:302:HOH:O	2.21	0.45
1:D:79:ALA:HA	2:D:201:EOH:H21	2.00	0.44
1:A:58:TYR:HB2	1:C:56:ARG:HG2	1.99	0.44
1:A:11:ASN:ND2	1:A:77:ASN:HB3	2.33	0.44
1:B:78:PRO:HG2	1:B:82:THR:OG1	2.18	0.44
1:A:96:ASN:HB2	1:C:21:ARG:NH1	2.33	0.43
1:D:1:MET:CE	1:D:2:THR:O	2.56	0.43
1:D:11:ASN:ND2	1:D:77:ASN:HB3	2.34	0.43
1:D:78:PRO:HG2	1:D:82[B]:THR:HB	1.99	0.43
1:C:78:PRO:HG2	1:C:82:THR:OG1	2.19	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116[A]:HIS:HE1	1:B:118:TYR:CE2	2.36	0.43
1:D:78:PRO:HG2	1:D:82[A]:THR:OG1	2.20	0.42
1:D:74:ILE:O	1:D:99:PHE:HA	2.19	0.42
1:B:47:VAL:O	4:B:303:HOH:O	2.22	0.42
1:C:58:TYR:CE1	1:C:59:GLU:HG3	2.54	0.42
1:A:78:PRO:HG2	1:A:82:THR:OG1	2.20	0.42
1:A:74:ILE:O	1:A:99:PHE:HA	2.19	0.41
1:A:96:ASN:HB2	1:C:21:ARG:HH11	1.85	0.41
1:A:96:ASN:HB2	1:C:21:ARG:HE	1.86	0.41
1:B:78:PRO:HG3	1:B:119:LEU:HD12	2.03	0.41
1:B:74:ILE:O	1:B:99:PHE:HA	2.21	0.41
1:D:44:LEU:HA	1:D:44:LEU:HD23	1.79	0.41
1:B:56:ARG:HG2	1:C:58:TYR:CB	2.50	0.40
1:C:18:LEU:HB2	4:C:303:HOH:O	2.20	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	147/152 (97%)	143 (97%)	4 (3%)	0	100	100
1	B	150/152 (99%)	145 (97%)	5 (3%)	0	100	100
1	C	147/152 (97%)	142 (97%)	4 (3%)	1 (1%)	18	14
1	D	151/152 (99%)	146 (97%)	4 (3%)	1 (1%)	18	14
All	All	595/608 (98%)	576 (97%)	17 (3%)	2 (0%)	36	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	2	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	-1	SER

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	126/128 (98%)	123 (98%)	3 (2%)	43	47
1	B	126/128 (98%)	119 (94%)	7 (6%)	19	16
1	C	124/128 (97%)	118 (95%)	6 (5%)	23	21
1	D	129/128 (101%)	121 (94%)	8 (6%)	16	13
All	All	505/512 (99%)	481 (95%)	24 (5%)	23	21

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	32	ASP
1	A	45	GLN
1	B	3	GLN
1	B	20	GLN
1	B	24	GLU
1	B	45	GLN
1	B	49	LEU
1	B	112	SER
1	B	148	ASN
1	C	20	GLN
1	C	21	ARG
1	C	24	GLU
1	C	28	SER
1	C	45	GLN
1	C	63	LYS
1	D	1	MET
1	D	3	GLN
1	D	20	GLN
1	D	24	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	32	ASP
1	D	36	LYS
1	D	45	GLN
1	D	148	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	115	HIS
1	B	11	ASN
1	B	20	GLN
1	B	115	HIS
1	B	140	GLN
1	C	11	ASN
1	C	115	HIS
1	D	3	GLN
1	D	4	GLN
1	D	11	ASN
1	D	115	HIS
1	D	140	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](#)

14 ligands 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	$\# Z > 2$
3	SO4	D	204	-	4,4,4	1.02	0	6,6,6	0.52	0
3	SO4	B	202	-	4,4,4	0.46	0	6,6,6	0.74	0
3	SO4	C	203	-	4,4,4	0.30	0	6,6,6	0.08	0
3	SO4	A	203	-	4,4,4	0.34	0	6,6,6	0.08	0
3	SO4	D	202	-	4,4,4	1.34	0	6,6,6	1.50	1 (16%)
3	SO4	B	203	-	4,4,4	0.35	0	6,6,6	0.45	0
3	SO4	D	203	-	4,4,4	0.79	0	6,6,6	0.31	0
3	SO4	A	202	-	4,4,4	0.50	0	6,6,6	0.66	0
2	EOH	A	201	-	2,2,2	0.48	0	1,1,1	0.15	0
2	EOH	C	201	-	2,2,2	0.20	0	1,1,1	0.40	0
3	SO4	B	204	-	4,4,4	0.51	0	6,6,6	0.21	0
2	EOH	B	201	-	2,2,2	0.65	0	1,1,1	0.13	0
2	EOH	D	201	-	2,2,2	0.05	0	1,1,1	0.77	0
3	SO4	C	202	-	4,4,4	0.40	0	6,6,6	0.60	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	202	SO4	O4-S-O2	3.16	126.08	109.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	202	SO4	1	0
3	B	203	SO4	1	0
2	B	201	EOH	1	0
2	D	201	EOH	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	148/152 (97%)	0.37	3 (2%) 65 64	48, 94, 147, 176	3 (2%)
1	B	151/152 (99%)	0.11	3 (1%) 65 64	40, 75, 130, 168	4 (2%)
1	C	149/152 (98%)	0.32	3 (2%) 65 64	59, 91, 152, 178	3 (2%)
1	D	152/152 (100%)	-0.29	1 (0%) 84 84	24, 56, 118, 174	2 (1%)
All	All	600/608 (98%)	0.12	10 (1%) 69 68	24, 82, 145, 178	12 (2%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-1	SER	3.9
1	A	116[A]	HIS	3.2
1	C	44	LEU	3.0
1	A	2	THR	2.9
1	B	0	HIS	2.9
1	D	149	ILE	2.6
1	B	44	LEU	2.5
1	A	132	ALA	2.3
1	C	146	LEU	2.2
1	C	149	ILE	2.1

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

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(Å ²)	Q<0.9
3	SO4	D	204	5/5	0.67	0.12	73,75,95,96	0
3	SO4	B	204	5/5	0.72	0.11	87,91,107,108	0
3	SO4	A	203	5/5	0.72	0.09	138,147,158,176	0
3	SO4	C	203	5/5	0.78	0.08	141,159,163,182	0
3	SO4	D	203	5/5	0.81	0.09	88,100,109,141	0
3	SO4	B	203	5/5	0.88	0.08	86,98,107,134	0
3	SO4	C	202	5/5	0.92	0.09	76,88,97,109	0
2	EOH	C	201	3/3	0.92	0.21	77,77,93,101	0
3	SO4	A	202	5/5	0.93	0.10	68,72,87,89	0
2	EOH	A	201	3/3	0.94	0.14	68,68,68,86	0
3	SO4	B	202	5/5	0.94	0.09	58,68,75,77	0
2	EOH	B	201	3/3	0.95	0.19	58,58,69,81	0
2	EOH	D	201	3/3	0.96	0.10	43,43,46,59	0
3	SO4	D	202	5/5	0.97	0.06	47,48,56,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.