



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:52 PM GMT

PDB ID : 4ATD
Title : Crystal structure of native Raucaffricine glucosidase
Authors : Xia, L.; Rajendran, C.; Ruppert, M.; Panjekar, S.; Wang, M.; Stoeckigt, J.
Deposited on : 2012-05-05
Resolution : 2.10 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

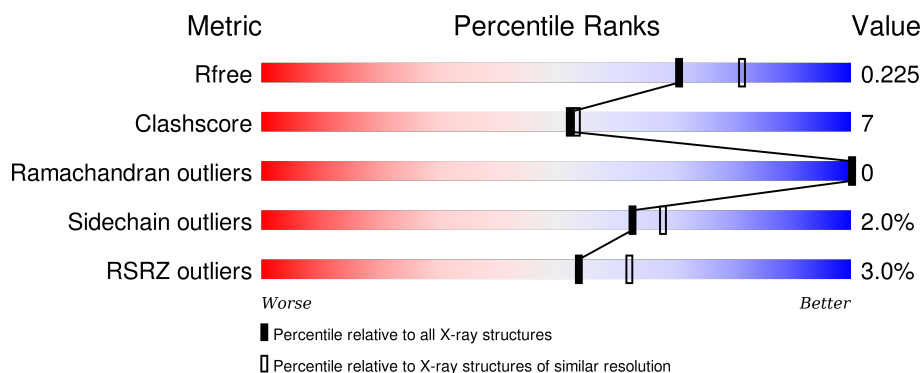
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	513	<div> <div>3%</div> <div>78%</div> <div>12%</div> <div>••</div> <div>9%</div> </div>
1	B	513	<div> <div>3%</div> <div>81%</div> <div>9%</div> <div>•</div> <div>9%</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
2	SO4	A	1515	-	-	-	X
2	SO4	A	1516	-	-	-	X
2	SO4	B	1515	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8111 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 RAUCAFFRICINE-O-BETA-D-GLUCOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	469	Total	C	N	O	S	0	4	0
			3793	2427	648	705	13			
1	B	469	Total	C	N	O	S	0	6	0
			3805	2436	649	707	13			

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

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

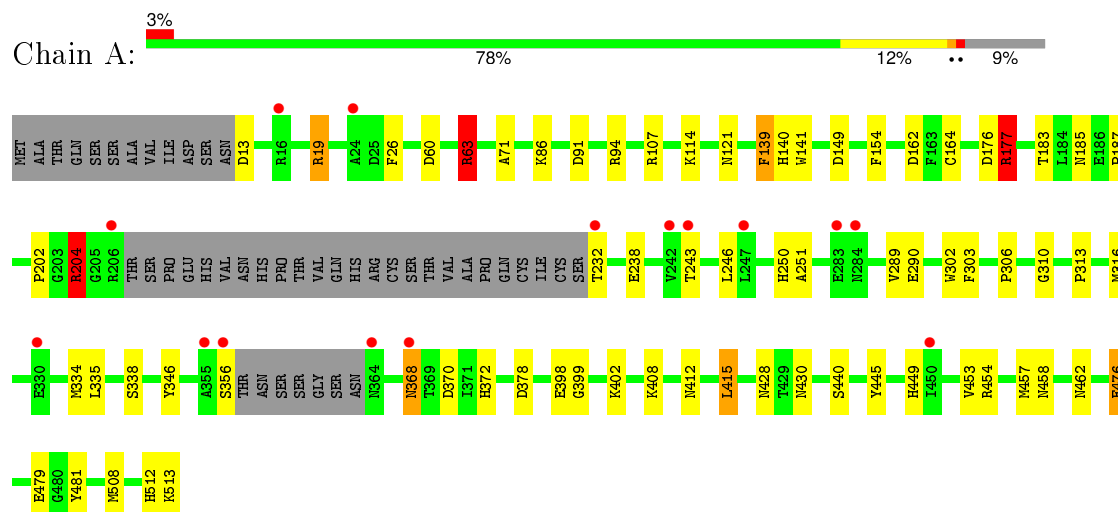
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	248	Total	O	0	0
			248	248		
3	B	240	Total	O	0	0
			240	240		

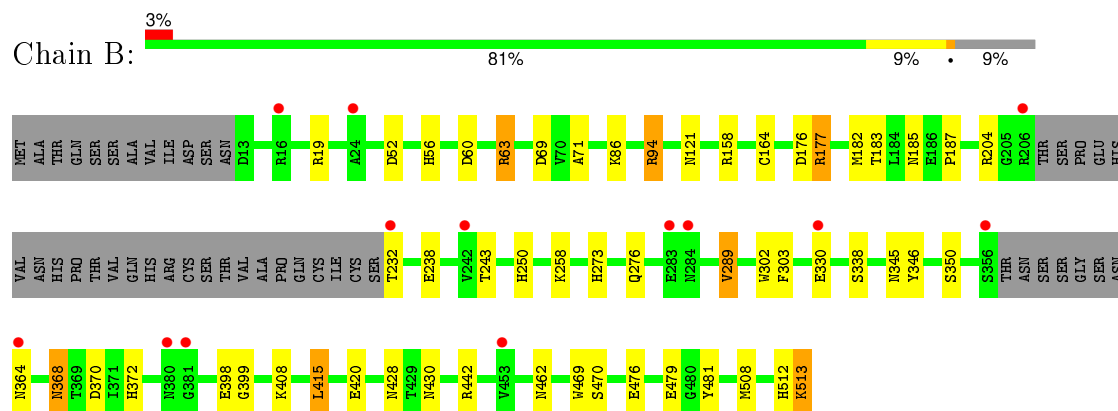
3 Residue-property plots

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 ($RSRZ > 2$). 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: RAUCAFFRICINE-O-BETA-D-GLUCOSIDASE



• Molecule 1: RAUCAFFRICINE-O-BETA-D-GLUCOSIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	102.24Å 127.25Å 216.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.95 – 2.10 19.95 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.95-2.10) 99.9 (19.95-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.85 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.175 , 0.206 0.193 , 0.225	Depositor DCC
R_{free} test set	977 reflections (1.20%)	DCC
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 82120 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8111	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.15	9/3916 (0.2%)	1.01	21/5313 (0.4%)
1	B	1.17	5/3934 (0.1%)	0.97	17/5336 (0.3%)
All	All	1.16	14/7850 (0.2%)	0.99	38/10649 (0.4%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	289	VAL	CB-CG1	-10.36	1.31	1.52
1	B	164	CYS	CB-SG	-7.28	1.69	1.82
1	B	289	VAL	CB-CG2	-6.59	1.39	1.52
1	A	289	VAL	CB-CG2	-6.55	1.39	1.52
1	A	289	VAL	CB-CG1	-6.46	1.39	1.52
1	A	164	CYS	CB-SG	-6.10	1.71	1.82
1	A	440	SER	CB-OG	-5.91	1.34	1.42
1	A	398	GLU	CG-CD	5.86	1.60	1.51
1	B	330	GLU	CG-CD	5.82	1.60	1.51
1	A	479	GLU	CD-OE1	5.81	1.32	1.25
1	A	476	GLU	CD-OE1	5.61	1.31	1.25
1	B	398	GLU	CG-CD	5.56	1.60	1.51
1	A	402	LYS	CE-NZ	-5.12	1.36	1.49
1	A	139	PHE	CE2-CZ	5.11	1.47	1.37

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	204	ARG	NE-CZ-NH2	-15.49	112.56	120.30
1	A	204	ARG	NE-CZ-NH1	13.29	126.94	120.30
1	B	94	ARG	NE-CZ-NH2	-9.86	115.37	120.30
1	A	94	ARG	NE-CZ-NH2	-9.24	115.68	120.30
1	A	63	ARG	NE-CZ-NH2	8.75	124.67	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	63	ARG	NE-CZ-NH1	-8.64	115.98	120.30
1	B	442	ARG	NE-CZ-NH1	7.89	124.25	120.30
1	A	177[A]	ARG	NE-CZ-NH2	-7.63	116.48	120.30
1	A	177[B]	ARG	NE-CZ-NH2	-7.63	116.48	120.30
1	A	13	ASP	CB-CG-OD1	7.39	124.95	118.30
1	A	177[A]	ARG	NE-CZ-NH1	7.11	123.86	120.30
1	A	177[B]	ARG	NE-CZ-NH1	7.11	123.86	120.30
1	B	442	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	B	60	ASP	CB-CG-OD1	6.59	124.23	118.30
1	A	63	ARG	NE-CZ-NH1	-6.41	117.09	120.30
1	B	94	ARG	NE-CZ-NH1	6.17	123.38	120.30
1	B	63	ARG	NE-CZ-NH2	6.15	123.37	120.30
1	B	177[A]	ARG	NE-CZ-NH1	6.11	123.36	120.30
1	B	177[B]	ARG	NE-CZ-NH1	6.11	123.36	120.30
1	A	94	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	B	177[A]	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	B	177[B]	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	B	232	THR	OG1-CB-CG2	-5.95	96.32	110.00
1	B	408	LYS	CD-CE-NZ	-5.85	98.25	111.70
1	A	378	ASP	CB-CG-OD1	5.65	123.39	118.30
1	A	290	GLU	OE1-CD-OE2	-5.64	116.53	123.30
1	A	204	ARG	CG-CD-NE	-5.62	100.00	111.80
1	A	289	VAL	CG1-CB-CG2	-5.50	102.10	110.90
1	A	408	LYS	CD-CE-NZ	-5.47	99.11	111.70
1	A	60	ASP	CB-CG-OD1	5.33	123.10	118.30
1	B	182	MET	CG-SD-CE	5.33	108.72	100.20
1	B	158	ARG	NE-CZ-NH1	-5.30	117.65	120.30
1	A	398	GLU	OE1-CD-OE2	-5.29	116.95	123.30
1	B	415	LEU	CA-CB-CG	5.24	127.34	115.30
1	A	162	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	415	LEU	CA-CB-CG	5.09	127.02	115.30
1	B	63	ARG	CG-CD-NE	-5.06	101.17	111.80
1	A	149	ASP	CB-CG-OD2	5.04	122.83	118.30

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	3793	0	3608	54	0
1	B	3805	0	3627	45	0
2	A	15	0	0	0	0
2	B	10	0	0	0	0
3	A	248	0	0	3	0
3	B	240	0	0	6	0
All	All	8111	0	7235	99	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 (99) 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:114:LYS:HE3	3:A:2087:HOH:O	1.38	1.19
1:A:121:ASN:ND2	1:A:177[A]:ARG:HH22	1.60	1.00
1:A:121:ASN:HD22	1:A:177[A]:ARG:HH22	1.10	0.99
1:B:121:ASN:ND2	1:B:177[A]:ARG:HH22	1.60	0.99
1:B:121:ASN:HD22	1:B:177[A]:ARG:HH22	1.12	0.93
1:A:63:ARG:HG2	1:A:63:ARG:HH11	1.36	0.90
1:B:258:LYS:HE3	3:B:2143:HOH:O	1.74	0.86
1:A:368:ASN:H	1:A:368:ASN:HD22	1.29	0.78
1:A:313:PRO:HD2	1:A:316:MET:HE2	1.65	0.78
1:B:512:HIS:O	1:B:513:LYS:HB2	1.80	0.78
1:B:364:ASN:HA	3:B:2183:HOH:O	1.85	0.76
1:A:428:ASN:OD1	1:A:430:ASN:HB2	1.89	0.73
1:A:453:VAL:HG12	1:A:457:MET:HE3	1.70	0.73
1:A:313:PRO:HD2	1:A:316:MET:CE	2.21	0.70
1:A:121:ASN:HD21	1:A:177[B]:ARG:HH22	1.38	0.69
1:B:176:ASP:OD2	1:B:177[A]:ARG:HD3	1.95	0.66
1:A:183:THR:OG1	1:A:250:HIS:HD2	1.78	0.66
1:A:302:TRP:HA	1:A:316:MET:HE1	1.76	0.66
1:B:508:MET:O	1:B:512:HIS:HB2	1.96	0.66
1:B:428:ASN:OD1	1:B:430:ASN:HB2	1.97	0.64
1:A:302:TRP:HA	1:A:316:MET:CE	2.26	0.64
1:A:176:ASP:OD2	1:A:177[A]:ARG:HD3	1.98	0.63
1:B:121:ASN:HD21	1:B:177[B]:ARG:HH22	1.45	0.63
1:B:52:ASP:O	1:B:56:HIS:HD2	1.81	0.63
1:A:63:ARG:CG	1:A:63:ARG:HH11	2.10	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:THR:OG1	1:B:250:HIS:HD2	1.81	0.62
1:A:346:TYR:OH	1:A:399:GLY:HA3	2.01	0.60
1:A:302:TRP:CA	1:A:316:MET:HE1	2.31	0.60
1:A:91:ASP:OD1	1:A:513:LYS:HE3	2.01	0.60
1:A:121:ASN:HD22	1:A:177[A]:ARG:NH2	1.93	0.59
1:B:185:ASN:HD21	1:B:345:ASN:HD21	1.51	0.58
1:A:508:MET:O	1:A:512:HIS:HB2	2.03	0.58
1:A:121:ASN:ND2	1:A:177[B]:ARG:HH22	2.00	0.58
1:B:121:ASN:HD22	1:B:177[A]:ARG:NH2	1.93	0.56
1:B:370:ASP:O	1:B:372:HIS:HD2	1.89	0.56
1:B:346:TYR:OH	1:B:399:GLY:HA3	2.06	0.56
1:B:121:ASN:HD22	1:B:177[B]:ARG:HH12	1.51	0.56
1:A:121:ASN:ND2	1:A:177[A]:ARG:NH2	2.43	0.55
1:A:372:HIS:HA	3:A:2183:HOH:O	2.06	0.55
1:B:476:GLU:HG3	1:B:476:GLU:O	2.05	0.55
1:A:19:ARG:NH1	1:A:513:LYS:HE2	2.22	0.55
1:A:453:VAL:HG12	1:A:457:MET:CE	2.37	0.54
1:A:415:LEU:HD12	1:A:462:ASN:ND2	2.22	0.54
1:A:63:ARG:NH1	1:A:63:ARG:HG2	2.17	0.53
1:A:187:PRO:HB2	1:A:303:PHE:CE2	2.44	0.53
1:B:512:HIS:O	1:B:513:LYS:CB	2.50	0.53
1:B:121:ASN:ND2	1:B:177[B]:ARG:HH12	2.08	0.52
1:B:121:ASN:ND2	1:B:177[B]:ARG:HH22	2.07	0.52
1:A:370:ASP:O	1:A:372:HIS:HD2	1.94	0.51
1:B:56:HIS:ND1	3:B:2041:HOH:O	2.35	0.49
1:A:243:THR:HG23	1:A:303:PHE:CE2	2.47	0.49
1:A:476:GLU:HG3	1:A:476:GLU:O	2.12	0.49
1:A:368:ASN:ND2	1:A:368:ASN:H	2.05	0.49
1:A:71:ALA:HA	1:A:481:TYR:OH	2.13	0.49
1:B:56:HIS:CE1	3:B:2041:HOH:O	2.66	0.48
1:A:251:ALA:HB2	1:A:335:LEU:HD23	1.95	0.48
1:B:415:LEU:HD12	1:B:462:ASN:ND2	2.28	0.48
1:B:121:ASN:ND2	1:B:177[A]:ARG:NH2	2.45	0.48
1:A:306:PRO:O	1:A:310:GLY:HA2	2.14	0.47
1:B:185:ASN:HD21	1:B:345:ASN:ND2	2.12	0.47
1:A:202:PRO:HG2	1:A:204:ARG:HG3	1.97	0.47
1:A:176:ASP:OD2	1:A:177[A]:ARG:CD	2.63	0.46
1:A:121:ASN:HD22	1:A:177[B]:ARG:HH12	1.62	0.46
1:A:415:LEU:HD11	3:A:2229:HOH:O	2.17	0.45
1:B:368:ASN:H	1:B:368:ASN:ND2	2.14	0.45
1:B:469:TRP:HA	1:B:470:SER:HA	1.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:ASN:ND2	1:B:345:ASN:HD21	2.14	0.45
1:A:140:HIS:CE1	1:A:185:ASN:ND2	2.85	0.45
1:B:187:PRO:HB2	1:B:303:PHE:CE1	2.52	0.44
1:B:52:ASP:O	1:B:56:HIS:CD2	2.68	0.44
1:A:19:ARG:CZ	1:A:513:LYS:HE2	2.49	0.43
1:A:368:ASN:N	1:A:368:ASN:HD22	2.06	0.43
1:A:454:ARG:NH1	1:A:458:ASN:OD1	2.52	0.43
1:A:121:ASN:ND2	1:A:177[B]:ARG:HH12	2.17	0.43
1:B:276:GLN:NE2	1:B:350:SER:HB2	2.34	0.43
1:A:302:TRP:HA	1:A:316:MET:HE3	2.00	0.42
1:A:26:PHE:CE1	1:A:457:MET:HE1	2.53	0.42
1:B:56:HIS:HE1	1:B:69:ASP:OD2	2.03	0.42
1:A:176:ASP:CG	1:A:177[A]:ARG:HD3	2.40	0.42
1:B:364:ASN:N	3:B:2183:HOH:O	2.53	0.42
1:A:445:TYR:CE2	1:A:449:HIS:CE1	3.07	0.42
1:B:273:HIS:CE1	1:B:303:PHE:HB3	2.55	0.42
1:B:250:HIS:CE1	1:B:338:SER:O	2.73	0.42
1:A:154:PHE:CE1	1:A:246:LEU:HD23	2.55	0.41
1:B:364:ASN:CA	3:B:2183:HOH:O	2.58	0.41
1:A:415:LEU:HD12	1:A:462:ASN:HD22	1.86	0.41
1:A:204:ARG:HD3	1:A:238:GLU:OE1	2.20	0.41
1:B:63:ARG:NH2	1:B:479:GLU:OE1	2.46	0.41
1:B:176:ASP:CG	1:B:177[A]:ARG:HD3	2.40	0.41
1:B:243:THR:HG21	1:B:302:TRP:CE2	2.55	0.41
1:A:250:HIS:CE1	1:A:338:SER:O	2.74	0.40
1:B:243:THR:HG21	1:B:302:TRP:NE1	2.36	0.40
1:A:313:PRO:CD	1:A:316:MET:HE2	2.42	0.40
1:B:204[A]:ARG:NH1	1:B:238:GLU:OE2	2.55	0.40
1:B:71:ALA:HA	1:B:481:TYR:OH	2.21	0.40
1:A:139:PHE:CZ	1:A:141:TRP:HA	2.56	0.40
1:B:469:TRP:CE2	1:B:470:SER:HB3	2.57	0.40
1:B:176:ASP:OD2	1:B:177[A]:ARG:CD	2.66	0.40
1:B:94:ARG:NH2	1:B:420:GLU:OE2	2.49	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	467/513 (91%)	449 (96%)	18 (4%)	0	100	100
1	B	469/513 (91%)	455 (97%)	14 (3%)	0	100	100
All	All	936/1026 (91%)	904 (97%)	32 (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	399/435 (92%)	387 (97%)	12 (3%)	48	51
1	B	401/435 (92%)	396 (99%)	5 (1%)	78	84
All	All	800/870 (92%)	783 (98%)	17 (2%)	63	66

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ARG
1	A	63	ARG
1	A	86	LYS
1	A	107	ARG
1	A	177[A]	ARG
1	A	177[B]	ARG
1	A	204	ARG
1	A	232	THR

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Mol	Chain	Res	Type
1	A	334	MET
1	A	356	SER
1	A	368	ASN
1	A	412	ASN
1	B	19	ARG
1	B	86	LYS
1	B	289	VAL
1	B	368	ASN
1	B	513	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	121	ASN
1	A	250	HIS
1	A	364	ASN
1	A	368	ASN
1	A	421	ASN
1	B	36	GLN
1	B	56	HIS
1	B	121	ASN
1	B	185	ASN
1	B	250	HIS
1	B	276	GLN
1	B	354	ASN
1	B	364	ASN
1	B	368	ASN
1	B	372	HIS
1	B	421	ASN
1	B	452	ASN
1	B	493	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

5 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	$\# Z > 2$
2	SO4	A	1514	-	4,4,4	0.42	0	6,6,6	0.91	1 (16%)
2	SO4	A	1515	-	4,4,4	0.90	0	6,6,6	2.23	2 (33%)
2	SO4	A	1516	-	4,4,4	0.76	0	6,6,6	0.79	0
2	SO4	B	1514	-	4,4,4	0.62	0	6,6,6	0.57	0
2	SO4	B	1515	-	4,4,4	0.80	0	6,6,6	0.78	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
2	SO4	A	1514	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1515	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1516	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1514	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1515	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1515	SO4	O2-S-O1	-4.51	95.20	109.50
2	A	1515	SO4	O4-S-O3	-2.62	98.31	108.98
2	A	1514	SO4	O2-S-O1	-2.07	102.92	109.50

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	469/513 (91%)	-0.24	15 (3%)	51 60	27, 36, 52, 66	0
1	B	469/513 (91%)	-0.24	13 (2%)	56 64	26, 34, 50, 67	0
All	All	938/1026 (91%)	-0.24	28 (2%)	54 62	26, 35, 52, 67	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	356	SER	5.0
1	B	284	ASN	4.1
1	B	380	ASN	4.0
1	A	284	ASN	4.0
1	B	356	SER	3.8
1	A	364	ASN	3.7
1	A	232	THR	3.3
1	B	364	ASN	3.2
1	B	381	GLY	3.2
1	A	206	ARG	2.9
1	B	16	ARG	2.9
1	A	247	LEU	2.6
1	A	330	GLU	2.6
1	B	283	GLU	2.6
1	A	24	ALA	2.6
1	B	206	ARG	2.6
1	B	330	GLU	2.5
1	A	16	ARG	2.4
1	A	450	ILE	2.4
1	B	242	VAL	2.4
1	B	453	VAL	2.4
1	A	243	THR	2.4
1	A	355	ALA	2.3
1	B	24	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	283	GLU	2.3
1	A	368	ASN	2.2
1	B	232	THR	2.0
1	A	242	VAL	2.0

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 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
2	SO4	A	1515	5/5	0.89	0.20	6.82	30,31,35,37	5
2	SO4	A	1516	5/5	0.95	0.19	6.10	29,33,37,37	5
2	SO4	B	1515	5/5	0.95	0.18	3.46	29,32,34,34	5
2	SO4	A	1514	5/5	0.93	0.16	-	28,30,35,39	5
2	SO4	B	1514	5/5	0.91	0.16	-	32,32,33,42	5

6.5 Other polymers [i](#)

There are no such residues in this entry.