



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:12 AM UTC

PDB ID : 1PGJ / pdb_00001pgj
Title : X-RAY STRUCTURE OF 6-PHOSPHOGLUCONATE DEHYDROGENASE
FROM THE PROTOZOAN PARASITE T. BRUCEI
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Deposited on : 1998-03-16
Resolution : 2.82 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

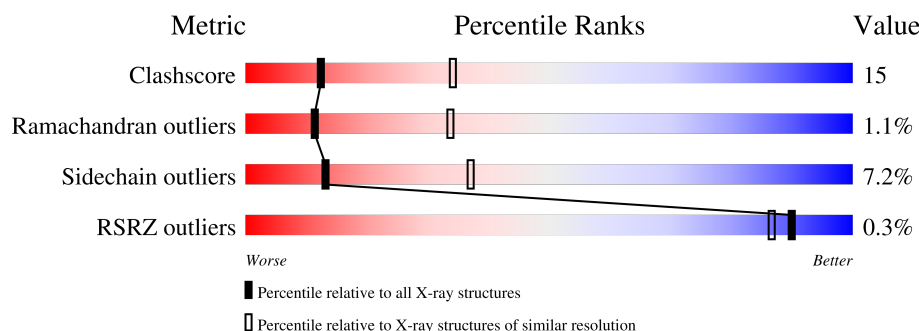
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	478	
1	B	478	

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	505	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7451 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 6-PHOSPHOGLUCONATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	478	Total	C	N	O	S	0	0	0
			3651	2309	629	688	25			
1	B	478	Total	C	N	O	S	0	0	0
			3651	2309	629	688	25			

- 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	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

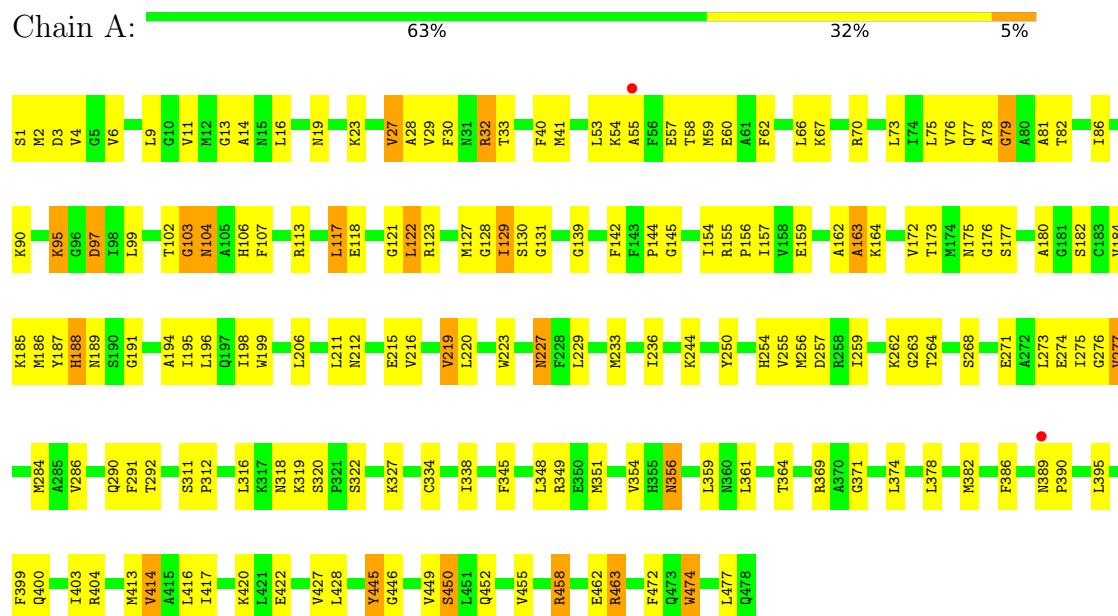
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	71	Total 71	O 71	0	0
3	B	58	Total 58	O 58	0	0

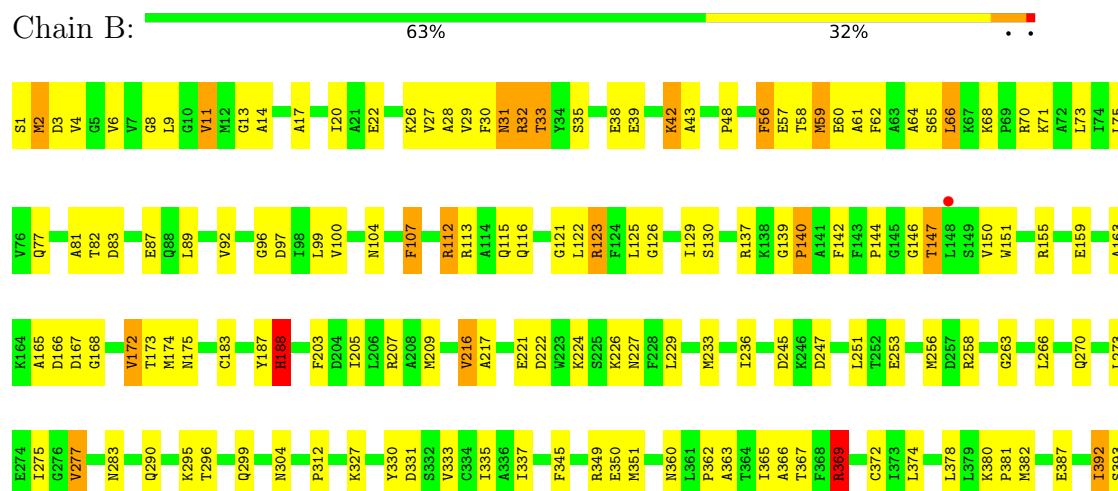
3 Residue-property plots

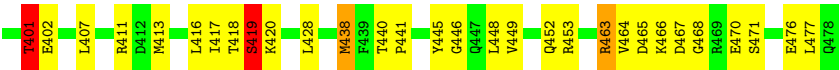
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 6-PHOSPHOGLUCONATE DEHYDROGENASE



• Molecule 1: 6-PHOSPHOGLUCONATE DEHYDROGENASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	135.13Å 135.13Å 116.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.76 – 2.82 19.76 – 2.82	Depositor EDS
% Data completeness (in resolution range)	97.7 (19.76-2.82) 97.4 (19.76-2.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.78 (at 2.83Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.186 , 0.273 0.200 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 74.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7451	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% 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: 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	0.77	0/3719	1.18	20/5015 (0.4%)
1	B	0.76	1/3719 (0.0%)	1.15	24/5015 (0.5%)
All	All	0.77	1/7438 (0.0%)	1.17	44/10030 (0.4%)

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	2
1	B	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	438	MET	SD-CE	-5.23	1.66	1.79

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	389	ASN	CA-C-N	10.71	133.22	119.84
1	A	389	ASN	C-N-CA	10.71	133.22	119.84
1	B	104	ASN	N-CA-C	9.83	125.03	111.17
1	B	66	LEU	N-CA-C	8.24	122.32	109.96
1	B	20	ILE	N-CA-C	7.38	117.47	110.53
1	A	369	ARG	N-CA-C	7.33	121.32	112.38
1	A	103	GLY	N-CA-C	7.29	130.46	113.18
1	A	227	ASN	N-CA-C	7.25	126.24	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	SER	N-CA-C	7.12	119.17	110.41
1	A	354	VAL	N-CA-C	7.11	117.21	110.53
1	B	56	PHE	N-CA-C	6.95	120.84	109.72
1	B	466	LYS	N-CA-C	-6.64	100.07	110.36
1	A	157	ILE	CB-CA-C	-6.50	103.64	111.97
1	B	476	GLU	N-CA-C	-6.14	101.84	110.50
1	A	117	LEU	N-CA-C	6.10	118.43	111.11
1	A	79	GLY	N-CA-C	-6.03	98.89	113.18
1	B	188	HIS	N-CA-C	-6.02	104.63	111.07
1	B	11	VAL	N-CA-C	-6.01	104.49	110.62
1	B	468	GLY	N-CA-C	-5.98	104.33	112.52
1	B	65	SER	N-CA-C	5.92	118.51	111.71
1	B	277	VAL	CB-CA-C	-5.70	105.64	110.94
1	B	175	ASN	N-CA-C	5.66	119.91	112.89
1	A	121	GLY	N-CA-C	5.65	121.81	115.08
1	A	188	HIS	N-CA-C	-5.57	105.11	111.07
1	B	126	GLY	N-CA-C	-5.56	100.01	113.18
1	A	356	ASN	N-CA-C	5.55	119.09	111.54
1	A	427	VAL	N-CA-C	5.46	116.23	110.72
1	A	474	TRP	N-CA-C	5.45	117.83	110.31
1	B	393	SER	N-CA-C	5.31	117.48	111.11
1	B	263	GLY	N-CA-C	5.30	125.74	113.18
1	B	27	VAL	N-CA-C	5.29	115.69	107.28
1	A	255	VAL	N-CA-C	5.28	116.31	108.23
1	B	419	SER	N-CA-C	5.28	117.11	111.36
1	B	369	ARG	N-CA-C	5.24	119.15	112.34
1	A	27	VAL	N-CA-C	5.21	115.35	107.75
1	B	227	ASN	N-CA-C	5.14	121.75	110.80
1	A	263	GLY	N-CA-C	5.08	125.23	113.18
1	A	277	VAL	CB-CA-C	-5.07	106.22	110.94
1	B	407	LEU	N-CA-C	5.06	117.19	111.11
1	B	60	GLU	N-CA-C	-5.06	105.66	111.07
1	B	392	ILE	N-CA-C	-5.03	102.36	109.21
1	B	216	VAL	N-CA-CB	5.03	118.13	110.58
1	A	163	ALA	N-CA-C	-5.03	103.04	110.48
1	B	387	GLU	N-CA-C	-5.03	105.51	111.69

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	187	TYR	Sidechain

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Mol	Chain	Res	Type	Group
1	A	445	TYR	Sidechain
1	B	187	TYR	Sidechain

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	3651	0	3655	114	0
1	B	3651	0	3655	125	0
2	A	10	0	0	2	0
2	B	10	0	0	0	0
3	A	71	0	0	4	0
3	B	58	0	0	4	0
All	All	7451	0	7310	225	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 15.

All (225) 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:66:LEU:HB2	1:A:70:ARG:HG3	1.47	0.95
1:B:13:GLY:HA2	1:B:75:LEU:HD13	1.56	0.88
1:B:142:PHE:HB2	1:B:172:VAL:HG13	1.55	0.88
1:A:57:GLU:HG2	1:A:58:THR:N	1.92	0.82
1:A:102:THR:HA	1:A:127:MET:O	1.79	0.81
1:B:32:ARG:HG3	1:B:33:THR:H	1.47	0.77
1:A:57:GLU:HG2	1:A:58:THR:H	1.47	0.77
1:A:139:GLY:O	1:A:163:ALA:HB2	1.84	0.77
1:B:233:MET:HE1	1:B:374:LEU:HD23	1.67	0.76
1:B:155:ARG:O	1:B:159:GLU:HG3	1.85	0.75
1:A:449:VAL:HA	1:A:452:GLN:OE1	1.89	0.73
1:A:446:GLY:O	1:A:450:SER:HB2	1.88	0.72
1:A:103:GLY:O	1:A:104:ASN:HB2	1.89	0.71
1:A:107:PHE:H	1:A:186:MET:HE1	1.54	0.71
1:A:41:MET:HE1	1:A:53:LEU:O	1.89	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:GLU:HG2	1:B:58:THR:N	2.06	0.70
1:B:125:LEU:HD13	1:B:150:VAL:CG2	2.21	0.69
1:A:378:LEU:HB3	1:A:382:MET:HE3	1.74	0.69
1:A:142:PHE:HB2	1:A:172:VAL:HG13	1.74	0.68
1:A:334:CYS:O	1:A:338:ILE:HG13	1.93	0.68
1:A:19:ASN:O	1:A:23:LYS:HG2	1.94	0.67
3:A:550:HOH:O	1:B:283:ASN:HB3	1.94	0.67
1:A:206:LEU:HD12	1:A:216:VAL:HG13	1.76	0.66
1:A:180:ALA:O	1:A:184:VAL:HG23	1.95	0.66
1:B:125:LEU:HD13	1:B:150:VAL:HG21	1.79	0.65
1:B:123:ARG:HB3	1:B:150:VAL:CG1	2.26	0.65
1:A:276:GLY:HA2	1:B:290:GLN:NE2	2.12	0.65
1:B:463:ARG:HD2	1:B:465:ASP:OD1	1.97	0.65
1:A:29:VAL:HG12	1:A:55:ALA:HA	1.78	0.64
1:A:312:PRO:HG2	1:A:420:LYS:HD3	1.78	0.64
1:A:322:SER:HA	1:A:327:LYS:HE3	1.79	0.64
1:B:70:ARG:HD3	1:B:97:ASP:OD2	1.98	0.63
1:B:30:PHE:CD1	1:B:31:ASN:N	2.66	0.63
1:A:188:HIS:HD2	1:A:189:ASN:HD22	1.47	0.63
1:B:378:LEU:O	1:B:381:PRO:HD2	1.99	0.63
1:B:6:VAL:HG13	1:B:75:LEU:HD12	1.80	0.62
1:B:378:LEU:O	1:B:382:MET:HG3	2.00	0.62
1:A:9:LEU:HD21	1:A:29:VAL:HG21	1.83	0.61
1:B:30:PHE:HB2	1:B:62:PHE:CD2	2.35	0.61
1:A:82:THR:O	1:A:86:ILE:HG13	1.99	0.61
1:B:73:LEU:HD23	1:B:100:VAL:HB	1.81	0.60
1:A:106:HIS:HA	1:A:186:MET:HE1	1.84	0.59
1:A:16:LEU:HD23	1:A:75:LEU:HD11	1.84	0.59
1:A:1:SER:C	1:A:2:MET:HE2	2.27	0.59
1:A:345:PHE:O	1:A:349:ARG:HG3	2.02	0.59
1:B:233:MET:CE	1:B:374:LEU:HD23	2.31	0.58
1:B:9:LEU:H	1:B:31:ASN:HD21	1.49	0.58
1:A:90:LYS:HE2	1:A:117:LEU:HD21	1.85	0.58
1:A:175:ASN:OD1	1:A:364:THR:HG23	2.05	0.57
1:A:212:ASN:OD1	1:A:215:GLU:HG3	2.05	0.57
1:A:102:THR:HB	1:A:129:ILE:HD11	1.87	0.57
1:B:57:GLU:HG2	1:B:58:THR:HG23	1.86	0.57
1:B:147:THR:O	1:B:150:VAL:HG22	2.05	0.56
1:A:70:ARG:HD3	1:A:97:ASP:OD2	2.05	0.56
1:A:196:LEU:HB3	1:B:452:GLN:HE22	1.71	0.56
1:B:245:ASP:HB3	1:B:247:ASP:OD1	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:ARG:HB3	1:A:156:PRO:HD3	1.87	0.56
1:A:458:ARG:NH2	1:A:472:PHE:O	2.39	0.55
1:A:122:LEU:N	1:A:122:LEU:HD23	2.20	0.55
1:A:477:LEU:HD13	1:B:236:ILE:HG22	1.89	0.55
1:A:131:GLY:HA2	1:A:371:GLY:O	2.07	0.55
1:A:275:ILE:HG13	1:A:277:VAL:HG23	1.88	0.54
1:B:411:ARG:NE	3:B:542:HOH:O	2.40	0.54
1:A:4:VAL:HG12	1:A:27:VAL:HG12	1.90	0.54
1:B:35:SER:O	1:B:38:GLU:HB2	2.07	0.54
1:A:2:MET:HE2	1:A:2:MET:N	2.23	0.54
1:A:142:PHE:HD2	1:A:172:VAL:HG22	1.73	0.54
1:A:2:MET:HG3	1:A:28:ALA:HB2	1.90	0.54
1:A:144:PRO:HD2	1:A:173:THR:O	2.08	0.54
1:A:191:GLY:O	1:A:195:ILE:HG12	2.08	0.53
1:A:199:TRP:HH2	1:A:229:LEU:HD13	1.72	0.53
1:B:57:GLU:CG	1:B:58:THR:HG23	2.38	0.53
1:A:220:LEU:N	1:A:220:LEU:HD12	2.24	0.53
1:B:256:MET:HG3	1:B:258:ARG:HD2	1.90	0.53
1:B:295:LYS:O	1:B:299:GLN:HG3	2.09	0.53
1:A:78:ALA:CB	1:A:81:ALA:HB3	2.40	0.52
1:A:474:TRP:HB2	3:A:569:HOH:O	2.10	0.52
1:B:59:MET:HG2	1:B:92:VAL:HG21	1.91	0.52
1:B:61:ALA:O	1:B:64:ALA:HB3	2.10	0.52
1:B:360:ASN:ND2	1:B:362:PRO:HG2	2.26	0.52
1:B:130:SER:O	1:B:140:PRO:HB2	2.09	0.51
1:A:445:TYR:O	1:A:449:VAL:HG23	2.10	0.51
1:A:14:ALA:HB1	1:A:40:PHE:CE1	2.45	0.51
1:A:162:ALA:HB2	1:A:172:VAL:CG2	2.41	0.51
1:B:96:GLY:O	1:B:123:ARG:NH1	2.43	0.51
1:B:107:PHE:CZ	1:B:351:MET:HE3	2.45	0.51
1:A:259:ILE:HD11	1:B:446:GLY:HA2	1.91	0.51
1:B:3:ASP:HB2	1:B:26:LYS:O	2.10	0.51
1:B:448:LEU:O	1:B:452:GLN:HG3	2.11	0.51
1:B:331:ASP:O	1:B:335:ILE:HG13	2.11	0.51
1:B:360:ASN:HD21	1:B:362:PRO:HG2	1.76	0.50
1:B:166:ASP:C	1:B:168:GLY:H	2.20	0.50
1:A:145:GLY:HA2	1:A:176:GLY:O	2.12	0.50
1:A:291:PHE:CD1	1:B:438:MET:HE3	2.46	0.50
1:B:57:GLU:CG	1:B:58:THR:N	2.75	0.50
1:A:130:SER:HA	3:A:555:HOH:O	2.12	0.50
1:A:107:PHE:N	1:A:186:MET:HE1	2.25	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:366:ALA:O	1:B:369:ARG:HD3	2.12	0.49
1:A:286:VAL:O	1:A:290:GLN:HG3	2.12	0.49
1:A:106:HIS:HA	1:A:186:MET:CE	2.42	0.49
1:A:128:GLY:O	1:A:142:PHE:HA	2.11	0.49
1:B:123:ARG:HB3	1:B:150:VAL:HG12	1.93	0.49
1:B:203:PHE:O	1:B:207:ARG:HG2	2.12	0.49
1:A:9:LEU:HD21	1:A:29:VAL:CG2	2.40	0.49
1:A:194:ALA:O	1:A:198:ILE:HG13	2.12	0.49
1:B:139:GLY:O	1:B:163:ALA:HB2	2.13	0.49
1:A:345:PHE:HD2	1:A:395:LEU:HD22	1.77	0.49
1:B:123:ARG:HB3	1:B:150:VAL:HG11	1.95	0.49
1:A:13:GLY:HA2	1:A:75:LEU:HD13	1.94	0.48
1:A:399:PHE:O	1:A:403:ILE:HG13	2.13	0.48
1:B:345:PHE:O	1:B:349:ARG:HG3	2.12	0.48
1:B:296:THR:HB	3:B:525:HOH:O	2.13	0.48
1:A:102:THR:HG22	1:A:127:MET:HE2	1.95	0.48
1:A:155:ARG:O	1:A:159:GLU:HG3	2.12	0.48
1:A:6:VAL:HG13	1:A:75:LEU:HD12	1.94	0.48
1:B:42:LYS:HG2	1:B:43:ALA:N	2.28	0.48
1:B:4:VAL:HG23	1:B:71:LYS:HB2	1.96	0.48
1:B:146:GLY:O	1:B:174:MET:HE3	2.13	0.48
1:A:30:PHE:HB2	1:A:62:PHE:CD2	2.49	0.48
1:B:57:GLU:HG2	1:B:58:THR:OG1	2.14	0.48
1:B:416:LEU:HD12	1:B:416:LEU:O	2.14	0.48
1:A:118:GLU:HB2	3:A:560:HOH:O	2.14	0.47
1:A:236:ILE:HG22	1:B:477:LEU:CD1	2.45	0.47
1:B:401:THR:HG22	1:B:402:GLU:N	2.28	0.47
1:A:414:VAL:HG22	1:B:418:THR:HG21	1.97	0.47
1:B:445:TYR:O	1:B:449:VAL:HG23	2.15	0.47
1:A:86:ILE:HD12	1:A:113:ARG:HD3	1.96	0.47
1:A:211:LEU:HD12	1:A:211:LEU:N	2.29	0.47
1:B:209:MET:HE1	1:B:413:MET:HE1	1.96	0.47
1:B:275:ILE:HG13	1:B:277:VAL:HG23	1.96	0.47
1:B:449:VAL:HA	1:B:452:GLN:NE2	2.30	0.46
1:B:449:VAL:HA	1:B:452:GLN:HE21	1.80	0.46
1:A:186:MET:HG2	1:A:268:SER:OG	2.14	0.46
1:B:428:LEU:HD23	1:B:428:LEU:HA	1.77	0.46
1:B:115:GLN:O	1:B:116:GLN:C	2.59	0.46
1:B:56:PHE:HD2	1:B:61:ALA:HB1	1.79	0.46
1:B:144:PRO:HD2	1:B:173:THR:O	2.16	0.46
1:B:150:VAL:HG23	1:B:151:TRP:N	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LEU:HD23	1:A:99:LEU:HA	1.78	0.46
1:B:89:LEU:HD13	1:B:99:LEU:HD22	1.97	0.46
1:B:62:PHE:CE1	1:B:66:LEU:HD11	2.52	0.46
1:B:205:ILE:HD13	1:B:417:ILE:CD1	2.46	0.46
1:B:188:HIS:CD2	1:B:188:HIS:C	2.94	0.45
1:B:440:THR:HA	1:B:441:PRO:HD3	1.85	0.45
1:B:68:LYS:HE2	1:B:70:ARG:NH2	2.31	0.45
1:B:277:VAL:HG22	1:B:350:GLU:HG3	1.97	0.45
1:A:361:LEU:HD23	1:A:386:PHE:CD1	2.51	0.45
1:B:83:ASP:O	1:B:87:GLU:HG3	2.17	0.45
1:B:9:LEU:N	1:B:31:ASN:HD21	2.13	0.45
1:B:39:GLU:C	1:B:42:LYS:HB3	2.42	0.44
1:A:103:GLY:O	1:A:104:ASN:CB	2.64	0.44
1:A:428:LEU:HD23	1:A:428:LEU:HA	1.91	0.44
1:A:102:THR:CB	1:A:129:ILE:HD11	2.48	0.44
1:B:9:LEU:HD21	1:B:29:VAL:HG13	2.00	0.44
1:B:363:ALA:O	1:B:367:THR:HG23	2.18	0.44
1:A:16:LEU:O	1:A:19:ASN:HB3	2.18	0.44
1:A:199:TRP:CH2	1:A:229:LEU:HD13	2.53	0.44
2:A:505:SO4:S	1:B:453:ARG:NH1	2.90	0.44
1:B:2:MET:HG2	1:B:28:ALA:HB2	1.99	0.44
1:B:327:LYS:O	1:B:330:TYR:HB3	2.18	0.44
1:A:233:MET:HE2	1:A:374:LEU:HD23	2.00	0.44
1:A:244:LYS:HE3	1:A:250:TYR:CE1	2.53	0.44
1:B:413:MET:O	1:B:417:ILE:HG13	2.18	0.43
1:A:463:ARG:HH22	1:B:245:ASP:CG	2.26	0.43
1:B:30:PHE:HE2	1:B:58:THR:CA	2.31	0.43
1:B:366:ALA:HA	1:B:369:ARG:HD3	2.01	0.43
1:A:14:ALA:HB1	1:A:40:PHE:HE1	1.82	0.43
1:B:59:MET:HB3	1:B:92:VAL:HG11	2.01	0.43
1:A:455:VAL:HG22	1:A:474:TRP:CE2	2.53	0.43
1:A:400:GLN:NE2	1:B:304:ASN:O	2.50	0.43
1:B:121:GLY:C	1:B:122:LEU:HD12	2.43	0.43
1:A:3:ASP:HA	1:A:67:LYS:HE2	2.00	0.43
1:A:254:HIS:CD2	1:A:254:HIS:N	2.87	0.43
1:B:14:ALA:O	1:B:17:ALA:HB3	2.19	0.43
1:B:327:LYS:O	1:B:327:LYS:HD2	2.19	0.43
1:A:4:VAL:O	1:A:27:VAL:HA	2.19	0.43
1:B:165:ALA:O	1:B:168:GLY:N	2.51	0.43
1:A:284:MET:HE2	3:B:549:HOH:O	2.18	0.42
1:A:291:PHE:CD1	1:B:438:MET:CE	3.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:GLN:O	1:B:81:ALA:HB3	2.19	0.42
1:B:205:ILE:HD13	1:B:417:ILE:HD13	2.01	0.42
1:A:271:GLU:HB3	1:A:351:MET:HE1	2.00	0.42
1:A:219:VAL:HG23	1:A:223:TRP:CE2	2.55	0.42
1:B:150:VAL:CG2	1:B:151:TRP:N	2.82	0.42
1:A:32:ARG:HD3	1:A:32:ARG:H	1.84	0.42
1:A:95:LYS:HD2	1:A:95:LYS:O	2.19	0.42
1:A:322:SER:CA	1:A:327:LYS:HE3	2.46	0.42
1:B:57:GLU:HG2	1:B:58:THR:CG2	2.49	0.42
1:B:256:MET:CG	1:B:258:ARG:HD2	2.49	0.42
1:A:27:VAL:O	1:A:27:VAL:HG23	2.19	0.42
1:A:462:GLU:HB2	1:B:256:MET:HB3	2.02	0.42
1:B:8:GLY:HA2	1:B:31:ASN:ND2	2.34	0.42
1:B:217:ALA:O	1:B:221:GLU:HB2	2.19	0.42
1:A:256:MET:HA	1:B:464:VAL:HG13	2.01	0.42
1:A:311:SER:HB2	1:A:312:PRO:HD2	2.01	0.42
1:A:400:GLN:O	1:A:404:ARG:HG3	2.20	0.42
1:B:183:CYS:SG	1:B:351:MET:HG2	2.60	0.42
1:B:362:PRO:HA	1:B:365:ILE:HG22	2.02	0.42
1:B:112:ARG:HG2	1:B:113:ARG:N	2.35	0.41
1:A:196:LEU:HB3	1:B:452:GLN:NE2	2.33	0.41
1:B:66:LEU:HB2	1:B:70:ARG:HG3	2.03	0.41
1:A:188:HIS:CD2	1:A:188:HIS:C	2.98	0.41
1:B:470:GLU:HG2	1:B:471:SER:N	2.36	0.41
1:A:257:ASP:O	1:A:292:THR:HB	2.20	0.41
1:A:316:LEU:O	1:A:319:LYS:NZ	2.54	0.41
1:B:365:ILE:O	1:B:365:ILE:HG12	2.20	0.41
1:B:251:LEU:HA	1:B:251:LEU:HD12	1.88	0.41
1:B:224:LYS:NZ	3:B:556:HOH:O	2.53	0.41
1:A:256:MET:HA	1:B:464:VAL:CG1	2.51	0.41
1:A:262:LYS:N	2:A:505:SO4:O2	2.54	0.41
1:B:82:THR:O	1:B:83:ASP:C	2.63	0.41
1:A:53:LEU:HA	1:A:53:LEU:HD23	1.89	0.41
1:B:22:GLU:CD	1:B:137:ARG:HH22	2.29	0.41
1:B:233:MET:HA	1:B:236:ILE:HG12	2.03	0.41
1:A:413:MET:O	1:A:417:ILE:HG13	2.21	0.40
1:B:57:GLU:HG2	1:B:58:THR:CB	2.51	0.40
1:B:266:LEU:O	1:B:270:GLN:HG3	2.21	0.40
1:B:419:SER:HB3	1:B:420:LYS:HD2	2.03	0.40
1:A:318:ASN:OD1	1:A:320:SER:HB3	2.22	0.40
1:B:32:ARG:HG3	1:B:33:THR:N	2.25	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:GLU:O	1:A:274:GLU:HB3	2.21	0.40
1:A:455:VAL:HG22	1:A:474:TRP:CD2	2.56	0.40
1:B:188:HIS:CE1	1:B:372:CYS:HB2	2.56	0.40
1:B:222:ASP:OD2	1:B:226:LYS:NZ	2.54	0.40
1:B:333:VAL:O	1:B:337:ILE:HG13	2.22	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	476/478 (100%)	445 (94%)	27 (6%)	4 (1%)	16	41
1	B	476/478 (100%)	439 (92%)	31 (6%)	6 (1%)	9	29
All	All	952/956 (100%)	884 (93%)	58 (6%)	10 (1%)	11	34

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	ASN
1	A	79	GLY
1	B	107	PHE
1	A	97	ASP
1	B	42	LYS
1	A	390	PRO
1	B	32	ARG
1	B	401	THR
1	B	140	PRO
1	B	48	PRO

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	381/381 (100%)	351 (92%)	30 (8%)	11	33
1	B	381/381 (100%)	356 (93%)	25 (7%)	15	41
All	All	762/762 (100%)	707 (93%)	55 (7%)	13	37

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	VAL
1	A	32	ARG
1	A	33	THR
1	A	54	LYS
1	A	59	MET
1	A	60	GLU
1	A	73	LEU
1	A	76	VAL
1	A	77	GLN
1	A	95	LYS
1	A	122	LEU
1	A	123	ARG
1	A	129	ILE
1	A	154	ILE
1	A	164	LYS
1	A	182	SER
1	A	185	LYS
1	A	219	VAL
1	A	227	ASN
1	A	264	THR
1	A	273	LEU
1	A	348	LEU
1	A	356	ASN
1	A	359	LEU
1	A	414	VAL
1	A	416	LEU
1	A	422	GLU

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Mol	Chain	Res	Type
1	A	450	SER
1	A	458	ARG
1	A	463	ARG
1	B	1	SER
1	B	2	MET
1	B	11	VAL
1	B	31	ASN
1	B	33	THR
1	B	59	MET
1	B	112	ARG
1	B	123	ARG
1	B	129	ILE
1	B	147	THR
1	B	167	ASP
1	B	172	VAL
1	B	188	HIS
1	B	216	VAL
1	B	229	LEU
1	B	253	GLU
1	B	273	LEU
1	B	312	PRO
1	B	369	ARG
1	B	380	LYS
1	B	392	ILE
1	B	401	THR
1	B	419	SER
1	B	463	ARG
1	B	467	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	ASN
1	A	106	HIS
1	A	188	HIS
1	A	283	ASN
1	B	15	ASN
1	B	31	ASN
1	B	116	GLN
1	B	188	HIS
1	B	213	ASN
1	B	283	ASN

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Mol	Chain	Res	Type
1	B	290	GLN
1	B	318	ASN
1	B	360	ASN
1	B	375	GLN
1	B	452	GLN
1	B	478	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](#)

4 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	505	-	4,4,4	0.70	0	6,6,6	0.65	0
2	SO4	B	505	-	4,4,4	0.99	0	6,6,6	0.43	0
2	SO4	B	507	-	4,4,4	0.74	0	6,6,6	0.78	0
2	SO4	A	507	-	4,4,4	0.55	0	6,6,6	0.91	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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	505	SO4	2	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	478/478 (100%)	-0.40	2 (0%) 88 84	4, 25, 51, 62	0
1	B	478/478 (100%)	-0.51	1 (0%) 91 89	5, 22, 48, 59	0
All	All	956/956 (100%)	-0.45	3 (0%) 90 86	4, 24, 50, 62	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	55	ALA	2.3
1	B	148	LEU	2.3
1	A	389	ASN	2.2

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
2	SO4	A	507	5/5	0.87	0.11	59,59,61,61	0
2	SO4	B	507	5/5	0.95	0.09	42,45,47,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	505	5/5	0.97	0.06	37,37,38,41	0
2	SO4	B	505	5/5	0.98	0.06	19,21,22,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.