



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:10 AM UTC

PDB ID : 1B7G / pdb_00001b7g
Title : GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE
Authors : Isupov, M.N.; Littlechild, J.A.
Deposited on : 1999-01-22
Resolution : 2.05 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

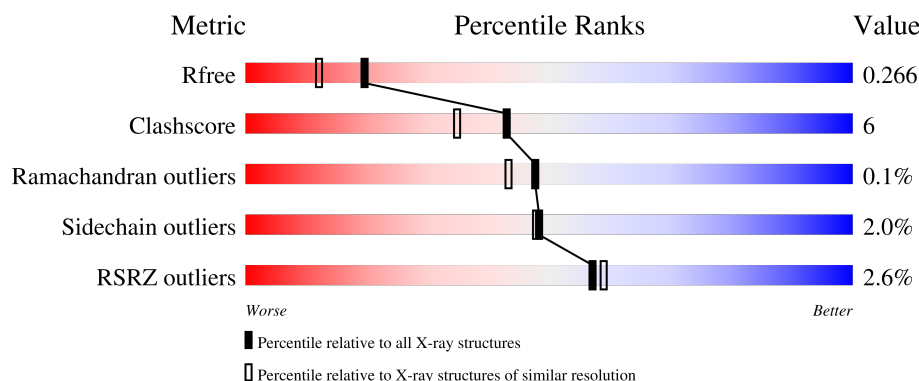
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	O	340	 84% 14% .
1	Q	340	 79% 19% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5844 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 (GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	O	340	Total	C	N	O	S	0	0	0
			2634	1664	456	500	14			
1	Q	340	Total	C	N	O	S	0	0	0
			2634	1664	456	500	14			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
O	2	VAL	ILE	SEE REMARK 999	UNP P39460
Q	2	VAL	ILE	SEE REMARK 999	UNP P39460
O	48	ILE	TYR	SEE REMARK 999	UNP P39460
Q	48	ILE	TYR	SEE REMARK 999	UNP P39460
O	?	-	ALA	deletion	UNP P39460
Q	?	-	ALA	deletion	UNP P39460
O	52	VAL	LEU	engineered mutation	UNP P39460
Q	52	VAL	LEU	engineered mutation	UNP P39460
O	72	GLU	ASP	SEE REMARK 999	UNP P39460
Q	72	GLU	ASP	SEE REMARK 999	UNP P39460
O	73	ASP	TYR	SEE REMARK 999	UNP P39460
Q	73	ASP	TYR	SEE REMARK 999	UNP P39460
O	159	LYS	ASN	SEE REMARK 999	UNP P39460
Q	159	LYS	ASN	SEE REMARK 999	UNP P39460

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



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

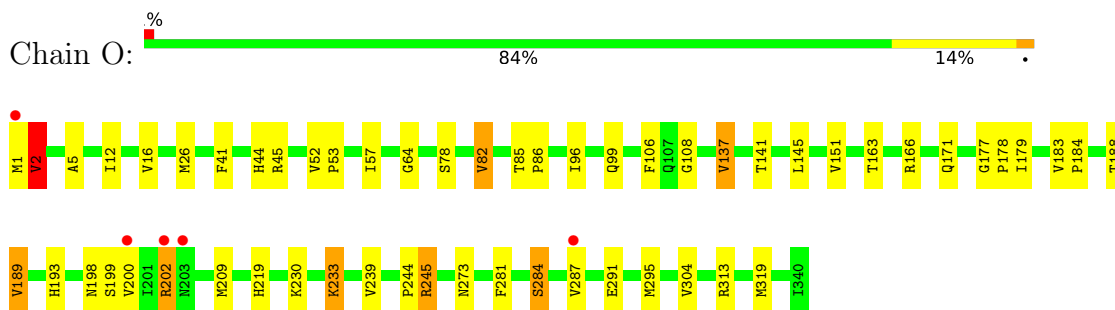
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	O	294	Total 294	O 294	0	0
3	Q	222	Total 222	O 222	0	0

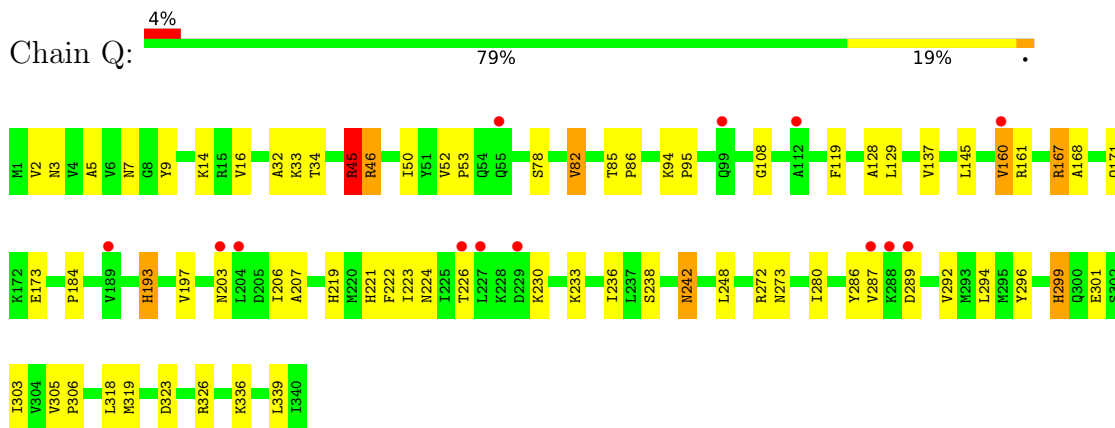
3 Residue-property plots [i](#)

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: PROTEIN (GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE)



- Molecule 1: PROTEIN (GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE)



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	101.55Å 101.55Å 179.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.00 – 2.05 22.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.6 (22.00-2.05) 99.5 (22.00-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 1.99Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.226 , 0.293 0.207 , 0.266	Depositor DCC
R_{free} test set	1298 reflections (2.03%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5844	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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	O	0.89	1/2670 (0.0%)	1.50	22/3613 (0.6%)
1	Q	0.86	1/2670 (0.0%)	1.56	29/3613 (0.8%)
All	All	0.87	2/5340 (0.0%)	1.53	51/7226 (0.7%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	Q	167	ARG	N-CA	5.26	1.52	1.46
1	O	284	SER	CA-C	5.09	1.59	1.52

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	119	PHE	N-CA-C	8.31	122.11	109.23
1	Q	242	ASN	CA-CB-CG	-8.06	104.53	112.60
1	Q	108	GLY	N-CA-C	7.85	123.30	114.67
1	Q	45	ARG	CD-NE-CZ	7.44	134.81	124.40
1	O	202	ARG	N-CA-C	7.14	119.06	111.28
1	Q	193	HIS	CA-CB-CG	-7.07	106.73	113.80
1	O	151	VAL	CA-C-N	7.01	130.61	120.38
1	O	151	VAL	C-N-CA	7.01	130.61	120.38
1	O	108	GLY	N-CA-C	6.90	122.06	114.40
1	Q	219	HIS	CA-CB-CG	6.82	120.61	113.80
1	O	106	PHE	CA-CB-CG	6.81	120.61	113.80
1	Q	160	VAL	CA-C-O	6.73	126.91	120.25
1	O	64	GLY	CA-C-N	-6.63	117.81	122.59
1	O	64	GLY	C-N-CA	-6.63	117.81	122.59
1	Q	119	PHE	CA-CB-CG	6.39	120.19	113.80
1	Q	272	ARG	CA-C-N	6.26	131.48	122.40
1	Q	272	ARG	C-N-CA	6.26	131.48	122.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	339	LEU	CA-C-O	6.25	127.14	120.70
1	Q	222	PHE	CA-CB-CG	-6.15	107.65	113.80
1	O	2	VAL	N-CA-CB	6.07	118.31	111.21
1	Q	336	LYS	CA-C-O	-6.04	114.99	121.45
1	Q	224	ASN	CA-C-N	-6.01	115.73	123.19
1	Q	224	ASN	C-N-CA	-6.01	115.73	123.19
1	Q	168	ALA	N-CA-C	5.81	117.28	111.07
1	Q	303	ILE	CA-C-O	-5.75	115.07	121.05
1	O	239	VAL	N-CA-CB	5.73	119.17	110.58
1	Q	82	VAL	N-CA-CB	5.71	118.33	111.31
1	O	193	HIS	CA-CB-CG	-5.61	108.19	113.80
1	O	184	PRO	CA-C-N	-5.55	116.86	122.74
1	O	184	PRO	C-N-CA	-5.55	116.86	122.74
1	O	189	VAL	O-C-N	5.53	127.41	121.10
1	O	188	THR	O-C-N	-5.53	116.71	122.96
1	Q	46	ARG	NE-CZ-NH1	-5.51	115.98	121.50
1	O	82	VAL	N-CA-CB	5.51	118.21	111.60
1	Q	203	ASN	CA-C-N	-5.50	114.45	122.65
1	Q	203	ASN	C-N-CA	-5.50	114.45	122.65
1	Q	184	PRO	CA-C-N	-5.43	116.99	122.74
1	Q	184	PRO	C-N-CA	-5.43	116.99	122.74
1	O	179	ILE	CB-CA-C	-5.42	104.73	112.22
1	Q	50	ILE	CA-C-O	5.33	126.23	120.53
1	O	281	PHE	CA-CB-CG	-5.29	108.51	113.80
1	O	202	ARG	CA-C-N	-5.22	114.38	122.67
1	O	202	ARG	C-N-CA	-5.22	114.38	122.67
1	Q	3	ASN	CA-CB-CG	-5.21	107.39	112.60
1	Q	160	VAL	N-CA-CB	5.11	119.55	111.99
1	O	44	HIS	CA-CB-CG	-5.11	108.69	113.80
1	Q	173	GLU	O-C-N	-5.07	116.69	122.93
1	Q	299	HIS	CA-C-N	5.06	127.57	120.28
1	Q	299	HIS	C-N-CA	5.06	127.57	120.28
1	O	199	SER	CA-C-N	5.05	128.18	123.08
1	O	199	SER	C-N-CA	5.05	128.18	123.08

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	O	2634	0	2739	28	0
1	Q	2634	0	2739	39	0
2	O	30	0	0	0	0
2	Q	30	0	0	0	0
3	O	294	0	0	6	0
3	Q	222	0	0	1	0
All	All	5844	0	5478	67	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 6.

All (67) 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:O:189:VAL:HG22	1:O:209:MET:HG2	1.68	0.76
1:Q:233:LYS:HA	1:Q:287:VAL:HG21	1.68	0.73
1:Q:230:LYS:HG3	1:Q:289:ASP:HA	1.72	0.71
1:O:198:ASN:O	1:O:202:ARG:HD3	1.91	0.69
1:Q:287:VAL:HG12	1:Q:292:VAL:HG22	1.74	0.68
1:O:291:GLU:HG3	3:O:742:HOH:O	1.97	0.64
1:Q:137:VAL:HG21	1:Q:145:LEU:HD12	1.80	0.64
1:Q:45:ARG:HH11	1:Q:45:ARG:HG3	1.68	0.59
1:O:244:PRO:O	1:O:245:ARG:HB2	2.04	0.58
1:O:2:VAL:HG12	1:O:26:MET:HG2	1.87	0.56
1:Q:167:ARG:HD2	1:Q:167:ARG:C	2.32	0.55
1:Q:5:ALA:HB2	1:Q:78:SER:OG	2.06	0.54
1:Q:85:THR:HB	1:Q:86:PRO:HD2	1.90	0.54
1:Q:236:ILE:HG13	1:Q:287:VAL:HG11	1.90	0.53
1:Q:233:LYS:HG3	1:Q:287:VAL:HG22	1.90	0.53
1:O:41:PHE:CE2	1:O:45:ARG:HD2	2.44	0.53
1:O:145:LEU:HD13	1:O:200:VAL:HG11	1.90	0.52
1:Q:193:HIS:O	1:Q:197:VAL:HG23	2.09	0.52
1:Q:160:VAL:O	1:Q:206:ILE:HA	2.11	0.51
1:Q:299:HIS:CE1	1:Q:301:GLU:HB2	2.46	0.50
1:Q:248:LEU:HD23	1:Q:280:ILE:HB	1.94	0.50
1:O:137:VAL:HB	1:O:141:THR:HB	1.94	0.50
1:O:52:VAL:HG13	1:O:53:PRO:HD2	1.93	0.49
1:O:233:LYS:HA	1:O:287:VAL:HG21	1.93	0.49
1:Q:94:LYS:N	1:Q:95:PRO:HD2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:171:GLN:HG3	3:Q:621:HOH:O	2.13	0.49
1:O:12:ILE:HG12	3:O:664:HOH:O	2.12	0.48
1:Q:161:ARG:HD3	1:Q:207:ALA:HB3	1.96	0.48
1:O:284:SER:CB	1:O:295:MET:HG2	2.43	0.47
1:O:313:ARG:HD3	1:O:319:MET:HE3	1.96	0.47
1:Q:9:TYR:CE1	1:Q:14:LYS:HA	2.50	0.47
1:Q:9:TYR:HB3	1:Q:33:LYS:HD2	1.96	0.47
1:Q:294:LEU:HD12	1:Q:294:LEU:C	2.39	0.47
1:Q:85:THR:HB	1:Q:86:PRO:CD	2.44	0.47
1:Q:323:ASP:OD1	1:Q:326:ARG:NH2	2.48	0.47
1:Q:52:VAL:HG12	1:Q:53:PRO:O	2.15	0.46
1:O:41:PHE:CZ	1:O:45:ARG:HD2	2.52	0.45
1:O:45:ARG:HB3	1:O:45:ARG:NH1	2.32	0.45
1:Q:16:VAL:HG11	1:Q:82:VAL:HG11	1.99	0.45
1:Q:223:ILE:HG22	1:Q:294:LEU:O	2.17	0.45
1:Q:34:THR:HG22	1:Q:53:PRO:HG2	1.99	0.44
1:Q:221:HIS:HB2	1:Q:296:TYR:CE1	2.53	0.44
1:Q:318:LEU:O	1:Q:319:MET:HB3	2.17	0.44
1:O:304:VAL:HG22	3:O:723:HOH:O	2.17	0.43
1:O:137:VAL:HG21	1:O:145:LEU:HD12	2.01	0.43
1:O:177:GLY:HA3	1:O:178:PRO:HA	1.75	0.43
1:Q:7:ASN:HA	1:Q:32:ALA:HB3	2.00	0.43
1:Q:46:ARG:HH11	1:Q:46:ARG:HD3	1.59	0.43
1:Q:2:VAL:HG21	1:Q:318:LEU:HD11	1.99	0.43
1:O:178:PRO:HG3	1:O:183:VAL:HG23	2.00	0.43
1:Q:305:VAL:HB	1:Q:306:PRO:HD3	1.99	0.42
1:O:96:ILE:HA	1:O:99:GLN:HG2	2.01	0.42
1:Q:128:ALA:O	1:Q:129:LEU:C	2.62	0.42
1:Q:161:ARG:HA	1:Q:207:ALA:O	2.20	0.42
1:Q:145:LEU:HD21	1:Q:197:VAL:HG13	2.02	0.42
1:O:202:ARG:HG2	3:O:720:HOH:O	2.20	0.42
1:O:85:THR:HB	1:O:86:PRO:CD	2.50	0.41
1:O:171:GLN:HG3	3:O:670:HOH:O	2.20	0.41
1:Q:160:VAL:HG13	1:Q:206:ILE:CG1	2.51	0.41
1:O:230:LYS:HG3	3:O:768:HOH:O	2.20	0.41
1:Q:287:VAL:HG12	1:Q:292:VAL:HG13	2.03	0.41
1:O:16:VAL:HG11	1:O:82:VAL:HG11	2.03	0.41
1:O:166:ARG:HA	1:O:219:HIS:ND1	2.37	0.40
1:Q:45:ARG:HH11	1:Q:45:ARG:CG	2.34	0.40
1:Q:286:TYR:CD1	1:Q:286:TYR:C	3.00	0.40
1:O:5:ALA:HB2	1:O:78:SER:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:244:PRO:O	1:O:245:ARG:CB	2.69	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	O	338/340 (99%)	328 (97%)	9 (3%)	1 (0%)	36	30
1	Q	338/340 (99%)	325 (96%)	13 (4%)	0	100	100
All	All	676/680 (99%)	653 (97%)	22 (3%)	1 (0%)	48	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	O	245	ARG

5.3.2 Protein sidechains [i](#)

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	O	295/295 (100%)	288 (98%)	7 (2%)	43	40
1	Q	295/295 (100%)	290 (98%)	5 (2%)	53	53
All	All	590/590 (100%)	578 (98%)	12 (2%)	48	48

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	1	MET
1	O	2	VAL
1	O	57	ILE
1	O	137	VAL
1	O	163	THR
1	O	233	LYS
1	O	273	ASN
1	Q	45	ARG
1	Q	226	THR
1	Q	238	SER
1	Q	242	ASN
1	Q	273	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	O	198	ASN
1	Q	87	ASN
1	Q	107	GLN
1	Q	198	ASN
1	Q	224	ASN
1	Q	308	ASN

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](#)

12 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	Q	612	-	4,4,4	0.65	0	6,6,6	0.24	0
2	SO4	Q	606	-	4,4,4	0.73	0	6,6,6	0.20	0
2	SO4	O	607	-	4,4,4	0.84	0	6,6,6	0.31	0
2	SO4	O	601	-	4,4,4	1.02	1 (25%)	6,6,6	0.81	0
2	SO4	O	603	-	4,4,4	0.68	0	6,6,6	0.48	0
2	SO4	O	605	-	4,4,4	0.69	0	6,6,6	0.09	0
2	SO4	Q	608	-	4,4,4	0.66	0	6,6,6	0.24	0
2	SO4	Q	602	-	4,4,4	0.71	0	6,6,6	0.31	0
2	SO4	O	610	-	4,4,4	0.70	0	6,6,6	0.07	0
2	SO4	Q	604	-	4,4,4	0.67	0	6,6,6	0.23	0
2	SO4	O	611	-	4,4,4	0.81	0	6,6,6	0.33	0
2	SO4	Q	609	-	4,4,4	0.71	0	6,6,6	0.25	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	O	601	SO4	O2-S	2.03	1.56	1.44

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

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 [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	O	340/340 (100%)	-0.15	5 (1%) 72 73	19, 30, 54, 87	0
1	Q	340/340 (100%)	0.44	13 (3%) 44 44	23, 37, 63, 85	0
All	All	680/680 (100%)	0.14	18 (2%) 57 59	19, 34, 59, 87	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	Q	287	VAL	3.7
1	Q	160	VAL	3.4
1	O	200	VAL	3.1
1	Q	227	LEU	2.6
1	Q	203	ASN	2.6
1	Q	55	GLN	2.5
1	Q	99	GLN	2.4
1	Q	288	LYS	2.4
1	Q	112	ALA	2.3
1	O	1	MET	2.2
1	Q	189	VAL	2.1
1	Q	204	LEU	2.1
1	Q	226	THR	2.1
1	O	202	ARG	2.1
1	O	287	VAL	2.1
1	O	203	ASN	2.0
1	Q	289	ASP	2.0
1	Q	229	ASP	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 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	O	610	5/5	0.72	0.15	80,80,83,84	0
2	SO4	Q	609	5/5	0.73	0.15	60,68,70,70	0
2	SO4	Q	606	5/5	0.86	0.15	58,60,61,64	0
2	SO4	O	611	5/5	0.86	0.19	53,60,65,66	0
2	SO4	Q	608	5/5	0.88	0.10	60,67,69,71	0
2	SO4	Q	612	5/5	0.88	0.20	59,59,62,63	5
2	SO4	O	607	5/5	0.89	0.22	36,51,53,55	0
2	SO4	O	605	5/5	0.92	0.07	65,66,70,70	0
2	SO4	Q	602	5/5	0.93	0.10	38,40,45,46	0
2	SO4	Q	604	5/5	0.94	0.14	50,54,58,60	0
2	SO4	O	603	5/5	0.96	0.11	44,53,56,57	0
2	SO4	O	601	5/5	0.97	0.06	25,29,30,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.