



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:16 PM UTC

PDB ID : 2PBG / pdb_00002pbg
Title : 6-PHOSPHO-BETA-D-GALACTOSIDASE FORM-B
Authors : Wiesmann, C.; Schulz, G.E.
Deposited on : 1997-02-21
Resolution : 2.50 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

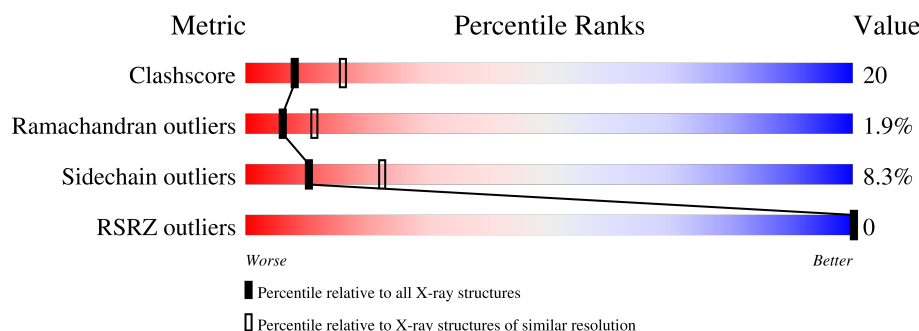
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	468	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4001 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-PHOSPHO-BETA-D-GALACTOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	468	3836	2475	624	726	11	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	256	CYS	SER	engineered mutation	UNP P11546

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



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

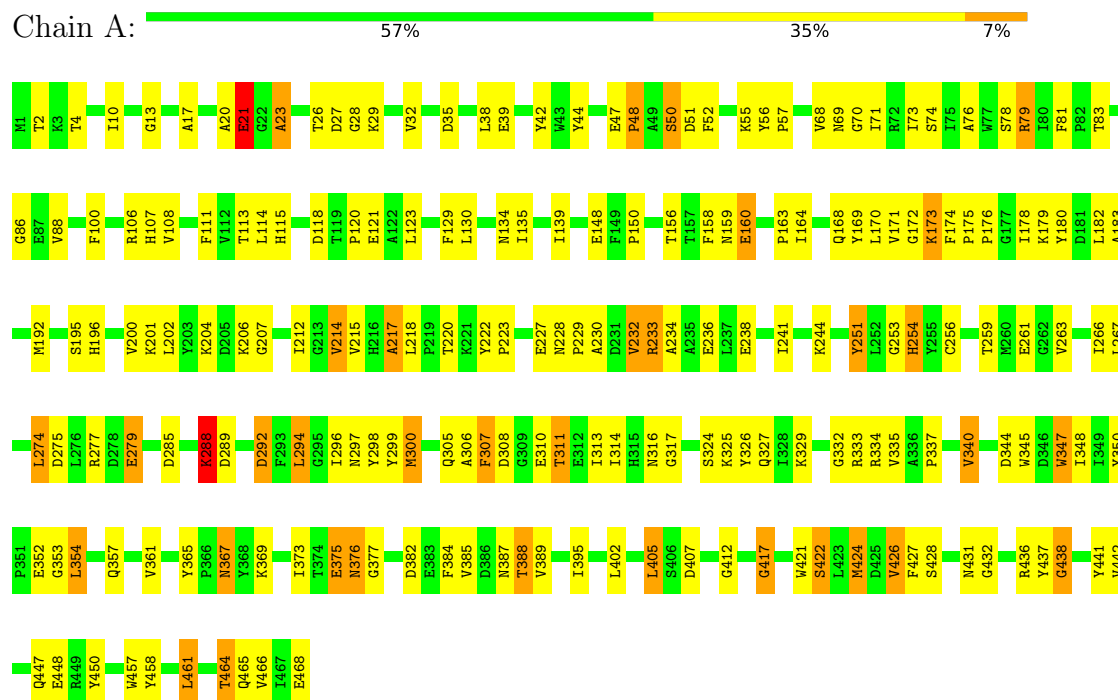
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	160	Total 160	O 160	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-PHOSPHO-BETA-D-GALACTOSIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	104.00Å 179.60Å 60.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.50 10.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.50) 76.8 (10.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.45Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.182 , 0.295 0.177 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	12.6	Xtriage
Anisotropy	0.762	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 74.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.024 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.034 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4001	wwPDB-VP
Average B, all atoms (Å ²)	10.0	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	0.56	0/3956	1.04	27/5367 (0.5%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	438	GLY	N-CA-C	9.63	124.87	112.54
1	A	27	ASP	N-CA-C	7.63	121.77	111.24
1	A	244	LYS	N-CA-C	7.23	119.16	111.28
1	A	130	LEU	N-CA-C	-6.91	104.04	113.30
1	A	412	GLY	N-CA-C	6.75	123.69	114.92
1	A	405	LEU	N-CA-C	-6.28	104.52	111.36
1	A	307	PHE	N-CA-C	6.26	118.51	107.61
1	A	340	VAL	N-CA-C	6.25	114.01	108.63
1	A	118	ASP	N-CA-C	6.10	119.91	111.55
1	A	365	TYR	CA-C-N	5.93	125.63	119.05
1	A	365	TYR	C-N-CA	5.93	125.63	119.05
1	A	28	GLY	N-CA-C	5.88	123.16	115.40
1	A	232	VAL	N-CA-C	-5.88	104.62	110.62
1	A	111	PHE	N-CA-C	-5.63	98.02	107.99
1	A	350	TYR	CA-C-N	5.47	124.93	119.24
1	A	350	TYR	C-N-CA	5.47	124.93	119.24
1	A	217	ALA	N-CA-C	-5.46	100.39	109.46
1	A	183	ALA	N-CA-C	-5.41	105.46	111.36
1	A	417	GLY	N-CA-C	5.40	118.37	110.38
1	A	160	GLU	N-CA-C	5.23	117.56	111.02
1	A	207	GLY	N-CA-C	5.22	120.90	114.69
1	A	347	TRP	N-CA-C	-5.11	100.92	109.24
1	A	236	GLU	N-CA-C	-5.10	105.41	110.97
1	A	388	THR	N-CA-C	-5.07	101.23	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	251	TYR	N-CA-C	5.03	120.26	113.72
1	A	288	LYS	N-CA-C	5.01	119.02	113.01
1	A	461	LEU	N-CA-C	-5.01	105.29	111.40

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	3836	0	3595	145	1
2	A	5	0	0	0	0
3	A	160	0	0	11	0
All	All	4001	0	3595	145	1

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 20.

All (145) 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:345:TRP:HE1	1:A:431:ASN:HD21	1.08	1.01
1:A:10:ILE:HD11	1:A:70:GLY:HA3	1.49	0.94
1:A:234:ALA:HB2	1:A:306:ALA:HB2	1.56	0.86
1:A:384:PHE:CZ	1:A:387:ASN:HA	2.16	0.81
1:A:345:TRP:HE1	1:A:431:ASN:ND2	1.80	0.79
1:A:267:LEU:HD21	1:A:274:LEU:HD13	1.67	0.77
1:A:307:PHE:H	1:A:332:GLY:HA2	1.49	0.77
1:A:316:ASN:ND2	1:A:345:TRP:HA	2.00	0.77
1:A:23:ALA:HB1	1:A:26:THR:HB	1.66	0.75
1:A:382:ASP:HB2	1:A:436:ARG:HD3	1.67	0.74
1:A:345:TRP:NE1	1:A:431:ASN:HD21	1.84	0.74
1:A:223:PRO:HB3	1:A:232:VAL:HA	1.71	0.72
1:A:424:MET:HE2	1:A:426:VAL:HG12	1.71	0.72
1:A:171:VAL:HG22	1:A:313:ILE:HG21	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:LEU:HD23	1:A:405:LEU:HD12	1.75	0.68
1:A:79:ARG:HG3	1:A:79:ARG:HH11	1.59	0.67
1:A:241:ILE:HD12	3:A:565:HOH:O	1.95	0.67
1:A:29:LYS:HD2	1:A:79:ARG:HH12	1.60	0.66
1:A:466:VAL:HG12	1:A:468:GLU:HG2	1.78	0.65
1:A:171:VAL:HG12	1:A:171:VAL:O	1.97	0.65
1:A:169:TYR:HA	1:A:178:ILE:HB	1.80	0.63
1:A:238:GLU:HA	3:A:565:HOH:O	1.98	0.63
1:A:388:THR:HG22	1:A:389:VAL:H	1.63	0.62
1:A:21:GLU:HA	1:A:52:PHE:HB3	1.81	0.62
1:A:251:TYR:HB3	1:A:367:ASN:HD21	1.65	0.61
1:A:86:GLY:HA3	3:A:596:HOH:O	1.99	0.60
1:A:228:ASN:O	1:A:232:VAL:HG23	2.01	0.60
1:A:253:GLY:O	1:A:254:HIS:HB3	2.01	0.60
1:A:447:GLN:HA	1:A:447:GLN:OE1	2.01	0.60
1:A:345:TRP:NE1	1:A:431:ASN:ND2	2.47	0.60
1:A:38:LEU:O	1:A:42:TYR:HB2	2.02	0.59
1:A:234:ALA:HB2	1:A:306:ALA:CB	2.29	0.59
1:A:305:GLN:OE1	1:A:335:VAL:HG21	2.02	0.58
1:A:428:SER:HB3	1:A:432:GLY:O	2.03	0.58
1:A:148:GLU:O	1:A:150:PRO:HD3	2.04	0.58
1:A:227:GLU:O	1:A:229:PRO:HD3	2.03	0.58
1:A:79:ARG:HG3	1:A:79:ARG:NH1	2.19	0.58
1:A:129:PHE:CD1	1:A:192:MET:HE3	2.39	0.58
1:A:314:ILE:HG12	1:A:325:LYS:HB3	1.86	0.57
1:A:10:ILE:HG13	1:A:69:ASN:OD1	2.05	0.57
1:A:202:LEU:HD11	1:A:206:LYS:NZ	2.19	0.57
1:A:56:TYR:CG	1:A:57:PRO:HD3	2.39	0.57
1:A:48:PRO:HG2	1:A:447:GLN:NE2	2.20	0.56
1:A:35:ASP:O	1:A:39:GLU:HG3	2.04	0.56
1:A:135:ILE:HG23	1:A:195:SER:OG	2.07	0.55
1:A:299:TYR:OH	1:A:375:GLU:HG2	2.07	0.55
1:A:68:VAL:O	1:A:108:VAL:HG22	2.06	0.54
1:A:38:LEU:HD22	1:A:427:PHE:HZ	1.73	0.54
1:A:426:VAL:HG23	1:A:427:PHE:O	2.07	0.54
1:A:316:ASN:HD21	1:A:345:TRP:HA	1.71	0.54
1:A:170:LEU:HD12	1:A:180:TYR:HA	1.90	0.54
1:A:222:TYR:OH	1:A:353:GLY:HA2	2.07	0.54
1:A:298:TYR:CD1	1:A:354:LEU:HG	2.43	0.53
1:A:382:ASP:OD1	1:A:436:ARG:HB2	2.09	0.53
1:A:168:GLN:HB3	1:A:176:PRO:HD2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:ALA:HA	1:A:50:SER:OG	2.09	0.52
1:A:200:VAL:HG13	1:A:212:ILE:HG21	1.90	0.52
1:A:377:GLY:HA3	3:A:504:HOH:O	2.10	0.52
1:A:327:GLN:HB3	3:A:585:HOH:O	2.11	0.51
1:A:395:ILE:HG23	1:A:457:TRP:CD2	2.46	0.51
1:A:357:GLN:O	1:A:361:VAL:HG23	2.11	0.50
1:A:17:ALA:O	1:A:21:GLU:HB2	2.11	0.50
1:A:29:LYS:HB2	1:A:79:ARG:NH1	2.27	0.50
1:A:71:ILE:HG22	1:A:108:VAL:HG12	1.92	0.50
1:A:218:LEU:HD12	1:A:296:ILE:HG12	1.93	0.50
1:A:10:ILE:HG23	1:A:417:GLY:HA3	1.92	0.50
1:A:384:PHE:CE2	1:A:387:ASN:HA	2.46	0.50
1:A:44:TYR:HD1	1:A:427:PHE:HE2	1.59	0.50
1:A:259:THR:O	1:A:263:VAL:HG23	2.12	0.50
1:A:337:PRO:O	1:A:340:VAL:HG12	2.12	0.50
1:A:196:HIS:O	1:A:200:VAL:HG23	2.11	0.49
1:A:233:ARG:O	1:A:233:ARG:HD3	2.11	0.49
1:A:442:VAL:HA	1:A:448:GLU:O	2.13	0.49
1:A:135:ILE:O	1:A:139:ILE:HD13	2.12	0.48
1:A:373:ILE:HG13	1:A:405:LEU:HD21	1.95	0.48
1:A:55:LYS:C	1:A:57:PRO:HD2	2.38	0.48
1:A:13:GLY:O	1:A:71:ILE:HA	2.14	0.48
1:A:285:ASP:O	1:A:288:LYS:HG3	2.13	0.48
1:A:81:PHE:HE1	1:A:88:VAL:HG22	1.78	0.48
1:A:29:LYS:HE3	3:A:491:HOH:O	2.14	0.48
1:A:182:LEU:HD22	1:A:266:ILE:CG2	2.44	0.48
1:A:81:PHE:CE1	1:A:88:VAL:HG22	2.49	0.47
1:A:297:ASN:OD1	1:A:375:GLU:HB2	2.14	0.47
1:A:196:HIS:CE1	1:A:200:VAL:HG21	2.49	0.47
1:A:175:PRO:HA	1:A:176:PRO:C	2.39	0.47
1:A:173:LYS:HD2	1:A:174:PHE:CE1	2.50	0.47
1:A:458:TYR:O	1:A:461:LEU:HB3	2.15	0.47
1:A:129:PHE:CG	1:A:192:MET:HG2	2.50	0.47
1:A:382:ASP:HB3	1:A:389:VAL:HG11	1.97	0.47
1:A:357:GLN:OE1	1:A:357:GLN:HA	2.15	0.47
1:A:106:ARG:O	1:A:107:HIS:HB2	2.16	0.46
1:A:56:TYR:N	1:A:57:PRO:CD	2.79	0.46
1:A:234:ALA:HB2	1:A:306:ALA:CA	2.46	0.46
1:A:448:GLU:HB3	1:A:450:TYR:CE1	2.51	0.46
1:A:51:ASP:HB3	1:A:55:LYS:HD3	1.98	0.46
1:A:171:VAL:O	1:A:171:VAL:CG1	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:GLU:HB2	3:A:527:HOH:O	2.16	0.46
1:A:106:ARG:HA	1:A:106:ARG:HD3	1.61	0.46
1:A:294:LEU:HD13	1:A:296:ILE:HG13	1.97	0.45
1:A:55:LYS:HD2	1:A:55:LYS:N	2.31	0.45
1:A:300:MET:HB2	1:A:348:ILE:O	2.16	0.45
1:A:212:ILE:O	1:A:292:ASP:HB2	2.15	0.45
1:A:305:GLN:HG2	1:A:307:PHE:HD2	1.82	0.45
1:A:317:GLY:O	1:A:431:ASN:ND2	2.49	0.45
1:A:44:TYR:CD1	1:A:427:PHE:CE2	3.05	0.45
1:A:156:THR:HG21	1:A:215:VAL:HG23	1.99	0.45
1:A:182:LEU:HD22	1:A:266:ILE:HG21	1.98	0.44
1:A:421:TRP:O	1:A:437:TYR:HB3	2.17	0.44
1:A:48:PRO:HG2	1:A:447:GLN:HE21	1.81	0.44
1:A:422:SER:O	1:A:438:GLY:HA2	2.17	0.44
1:A:464:THR:O	1:A:466:VAL:HG23	2.17	0.44
1:A:279:GLU:CD	1:A:279:GLU:H	2.25	0.44
1:A:23:ALA:CB	1:A:26:THR:HB	2.42	0.44
1:A:123:LEU:HD22	1:A:134:ASN:OD1	2.18	0.44
1:A:200:VAL:HG13	1:A:212:ILE:CG2	2.49	0.43
1:A:204:LYS:HD2	3:A:549:HOH:O	2.18	0.43
1:A:160:GLU:O	1:A:164:ILE:HG13	2.18	0.43
1:A:202:LEU:HD11	1:A:206:LYS:HZ2	1.82	0.43
1:A:47:GLU:OE1	1:A:48:PRO:HA	2.19	0.43
1:A:115:HIS:HB3	1:A:158:PHE:HD1	1.84	0.43
1:A:310:GLU:HB3	3:A:559:HOH:O	2.18	0.43
1:A:441:TYR:CD1	1:A:441:TYR:C	2.97	0.43
1:A:115:HIS:HB3	1:A:158:PHE:CD1	2.54	0.43
1:A:76:ALA:HB3	1:A:79:ARG:HB2	2.00	0.42
1:A:32:VAL:HG22	1:A:35:ASP:OD2	2.19	0.42
1:A:169:TYR:HD1	1:A:178:ILE:HD12	1.85	0.42
1:A:326:TYR:CD1	1:A:334:ARG:HG3	2.55	0.42
1:A:407:ASP:HA	3:A:608:HOH:O	2.19	0.42
1:A:120:PRO:O	1:A:121:GLU:C	2.63	0.42
1:A:123:LEU:HA	1:A:123:LEU:HD23	1.73	0.42
1:A:297:ASN:CG	1:A:375:GLU:HB2	2.45	0.41
1:A:275:ASP:OD1	1:A:277:ARG:HD2	2.20	0.41
1:A:375:GLU:O	1:A:376:ASN:HB2	2.20	0.41
1:A:311:THR:HB	1:A:329:LYS:H	1.85	0.41
1:A:402:LEU:HA	1:A:405:LEU:HD12	2.01	0.41
1:A:300:MET:HE3	1:A:347:TRP:HA	2.02	0.41
1:A:305:GLN:HG2	1:A:307:PHE:CD2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ILE:HD12	1:A:100:PHE:HE1	1.85	0.41
1:A:196:HIS:CE1	1:A:214:VAL:HG22	2.56	0.40
1:A:178:ILE:HG22	1:A:179:LYS:N	2.36	0.40
1:A:74:SER:HA	1:A:113:THR:O	2.21	0.40
1:A:160:GLU:HG2	1:A:217:ALA:HB2	2.04	0.40
1:A:230:ALA:O	1:A:306:ALA:HB2	2.22	0.40
1:A:333:ARG:HB2	3:A:585:HOH:O	2.21	0.40
1:A:344:ASP:HB2	1:A:345:TRP:CZ3	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:CYS:SG	1:A:256:CYS:SG[3_655]	2.19	0.01

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	466/468 (100%)	407 (87%)	50 (11%)	9 (2%)	6 11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	308	ASP
1	A	172	GLY
1	A	254	HIS
1	A	367	ASN
1	A	173	LYS
1	A	21	GLU
1	A	23	ALA
1	A	369	LYS

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Mol	Chain	Res	Type
1	A	376	ASN

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	397/397 (100%)	364 (92%)	33 (8%)	10	22

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	4	THR
1	A	21	GLU
1	A	48	PRO
1	A	50	SER
1	A	78	SER
1	A	79	ARG
1	A	83	THR
1	A	114	LEU
1	A	159	ASN
1	A	163	PRO
1	A	201	LYS
1	A	214	VAL
1	A	220	THR
1	A	233	ARG
1	A	261	GLU
1	A	274	LEU
1	A	279	GLU
1	A	288	LYS
1	A	289	ASP
1	A	292	ASP
1	A	294	LEU
1	A	300	MET
1	A	311	THR
1	A	324	SER

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Mol	Chain	Res	Type
1	A	354	LEU
1	A	375	GLU
1	A	385	VAL
1	A	422	SER
1	A	424	MET
1	A	426	VAL
1	A	464	THR
1	A	465	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	54	HIS
1	A	107	HIS
1	A	124	HIS
1	A	159	ASN
1	A	264	ASN
1	A	282	GLN
1	A	291	ASN
1	A	414	ASN
1	A	431	ASN
1	A	465	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](#)

1 ligand is 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	469	-	4,4,4	0.36	0	6,6,6	0.32	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.

No monomer is involved in short contacts.

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	468/468 (100%)	-0.68	0 100 100	2, 7, 26, 49	0

There are no RSRZ outliers to report.

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	469	5/5	0.92	0.09	39,40,42,43	5

6.5 Other polymers [i](#)

There are no such residues in this entry.