



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:54 AM UTC

PDB ID : 4AI8 / pdb_00004ai8
Title : FACTOR INHIBITING HIF-1 ALPHA IN COMPLEX WITH DAMINOZIDE
Authors : King, O.N.F.; Chowdhury, R.; Rose, N.R.; McDonough, M.A.; Clifton, I.J.; Schofield, C.J.; Kawamura, A.
Deposited on : 2012-02-08
Resolution : 2.40 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

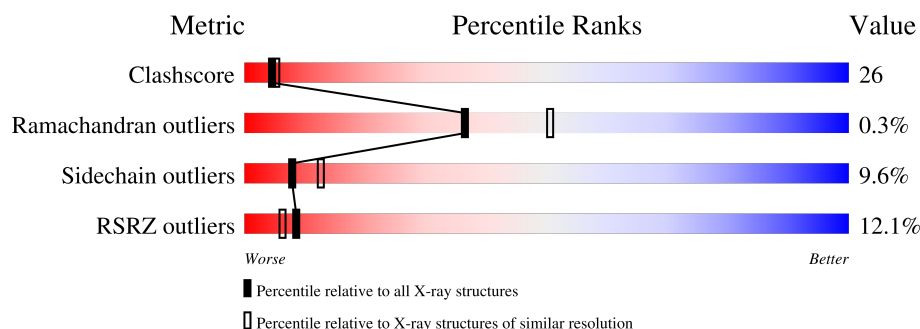
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	352	

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	1353	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 2897 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 HYPOXIA-INDUCIBLE FACTOR 1-ALPHA INHIBITOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	339	2707	1738	458	500	11	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q9NWT6
A	-1	SER	-	expression tag	UNP Q9NWT6
A	0	HIS	-	expression tag	UNP Q9NWT6

- 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		

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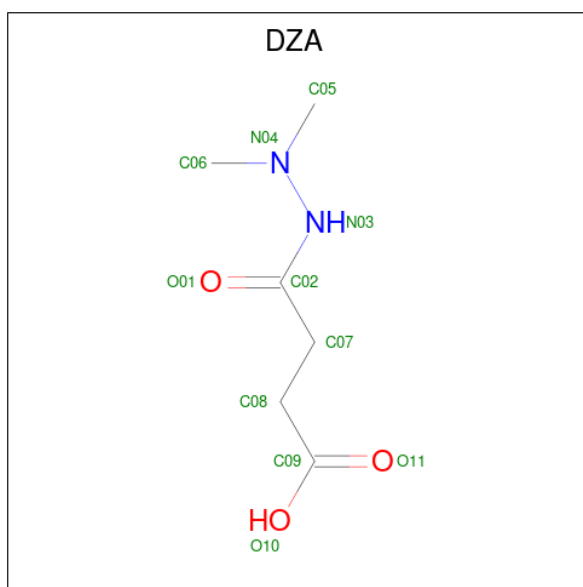
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is DAMINOZIDE (CCD ID: DZA) (formula: $C_6H_{12}N_2O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			11	6	2	3		

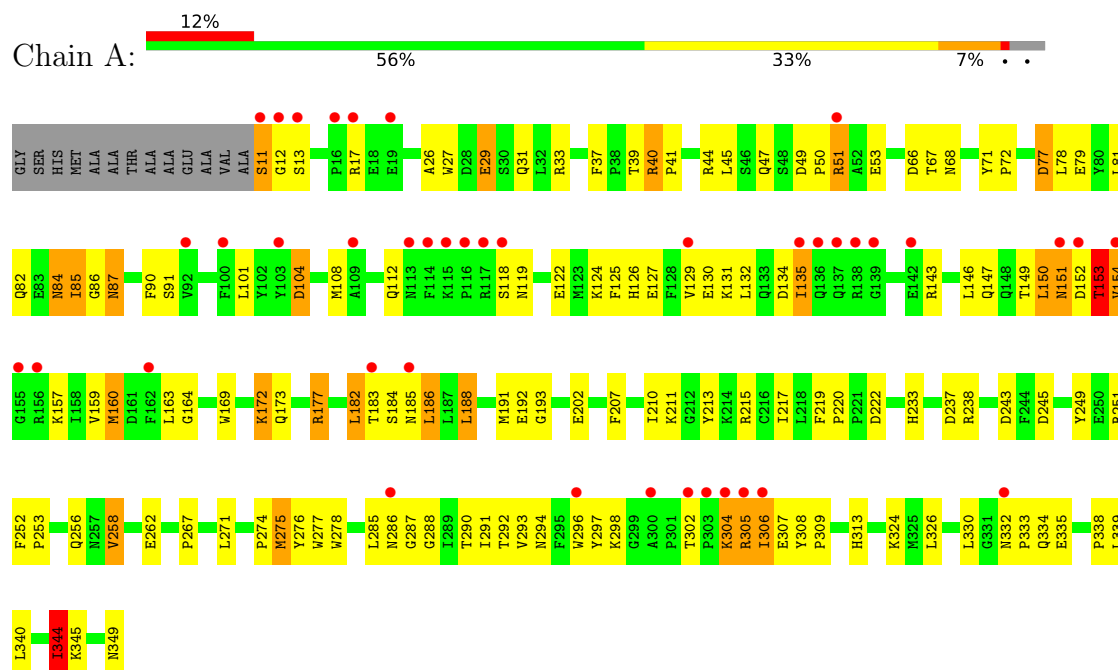
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	135	Total	O	0	0
			135	135		

- Molecule 1: HYPOXIA-INDUCIBLE FACTOR 1-ALPHA INHIBITOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	86.76Å 86.76Å 145.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.37 – 2.40 28.37 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.8 (28.37-2.40) 83.9 (28.37-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.39Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.231 , 0.248 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	48.5	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2897	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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: GOL, ZN, SO4, DZA

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	0/2789	1.02	14/3797 (0.4%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	84	ASN	N-CA-C	7.22	123.11	113.72
1	A	153	THR	N-CA-C	-6.45	103.70	113.89
1	A	298	LYS	N-CA-C	-6.39	100.82	110.28
1	A	66	ASP	N-CA-C	6.06	119.92	111.56
1	A	252	PHE	CA-C-N	5.58	125.68	119.32
1	A	252	PHE	C-N-CA	5.58	125.68	119.32
1	A	37	PHE	N-CA-C	5.53	117.64	110.40
1	A	154	VAL	N-CA-C	5.51	116.06	110.05
1	A	220	PRO	CA-C-N	5.24	125.30	119.32
1	A	220	PRO	C-N-CA	5.24	125.30	119.32
1	A	344	ILE	N-CA-C	5.24	117.13	111.58
1	A	77	ASP	N-CA-C	-5.22	101.33	108.38
1	A	258	VAL	N-CA-C	5.21	115.77	108.89
1	A	345	LYS	N-CA-C	5.17	117.66	109.96

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	2707	0	2534	139	1
2	A	25	0	0	3	0
3	A	18	0	24	0	0
4	A	11	0	11	0	0
5	A	1	0	0	0	0
6	A	135	0	0	5	0
All	All	2897	0	2569	139	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 26.

All (139) 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:332:ASN:HD21	1:A:334:GLN:HB2	1.13	1.14
1:A:125:PHE:O	1:A:129:VAL:HG23	1.61	1.01
1:A:160:MET:HE2	1:A:163:LEU:HD12	1.43	0.99
1:A:51:ARG:HG2	1:A:51:ARG:HH11	1.26	0.96
1:A:84:ASN:OD1	1:A:157:LYS:HD2	1.67	0.94
1:A:304:LYS:HG3	1:A:305:ARG:N	1.87	0.90
1:A:182:LEU:C	1:A:182:LEU:HD12	1.97	0.90
1:A:81:LEU:O	1:A:85:ILE:HG22	1.75	0.86
1:A:126:HIS:O	1:A:130:GLU:HG3	1.75	0.86
1:A:51:ARG:HH11	1:A:51:ARG:CG	1.90	0.85
1:A:51:ARG:HG2	1:A:51:ARG:NH1	1.91	0.82
1:A:304:LYS:HG3	1:A:305:ARG:H	1.42	0.82
1:A:151:ASN:ND2	1:A:153:THR:CG2	2.44	0.80
1:A:160:MET:HE2	1:A:160:MET:HA	1.63	0.80
1:A:31:GLN:O	1:A:285:LEU:HD12	1.80	0.80
1:A:173:GLN:OE1	1:A:177:ARG:NH1	2.13	0.80
1:A:151:ASN:HD21	1:A:153:THR:CG2	1.95	0.78
1:A:53:GLU:OE1	1:A:177:ARG:NH1	2.16	0.77
1:A:326:LEU:HD11	1:A:339:LEU:HD23	1.69	0.74
1:A:332:ASN:HD21	1:A:334:GLN:CB	1.98	0.73
1:A:151:ASN:HD21	1:A:153:THR:HG21	1.53	0.73
1:A:183:THR:HB	1:A:296:TRP:CD1	2.24	0.72
1:A:183:THR:CG2	1:A:296:TRP:HE1	2.03	0.71
1:A:147:GLN:HG2	1:A:188:LEU:HD23	1.72	0.71
1:A:91:SER:HA	1:A:122:GLU:OE1	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:ASN:ND2	1:A:334:GLN:HB2	1.98	0.70
1:A:172:LYS:HB3	6:A:2063:HOH:O	1.93	0.68
1:A:160:MET:HA	1:A:160:MET:CE	2.23	0.68
1:A:29:GLU:OE1	1:A:29:GLU:N	2.26	0.68
1:A:124:LYS:HB2	1:A:127:GLU:HG3	1.75	0.67
1:A:215:ARG:NE	1:A:262:GLU:OE2	2.27	0.67
1:A:33:ARG:NH2	1:A:193:GLY:O	2.26	0.67
1:A:151:ASN:ND2	1:A:153:THR:HG22	2.10	0.66
1:A:71:TYR:HB3	1:A:72:PRO:HD3	1.78	0.65
1:A:192:GLU:OE1	1:A:286:ASN:C	2.39	0.65
1:A:183:THR:HG21	1:A:296:TRP:HE1	1.59	0.65
1:A:182:LEU:HD12	1:A:184:SER:N	2.13	0.64
1:A:182:LEU:C	1:A:182:LEU:CD1	2.70	0.64
1:A:202:GLU:O	1:A:275:MET:CE	2.45	0.63
1:A:39:THR:O	1:A:40:ARG:HD3	1.98	0.63
1:A:151:ASN:CG	1:A:153:THR:HG22	2.24	0.63
1:A:219:PHE:HB2	1:A:278:TRP:HB2	1.80	0.62
1:A:154:VAL:HG23	1:A:159:VAL:HG23	1.82	0.62
1:A:210:ILE:CD1	1:A:293:VAL:HG23	2.30	0.62
1:A:211:LYS:CD	1:A:291:ILE:HD12	2.29	0.62
1:A:160:MET:CE	1:A:160:MET:CA	2.79	0.61
1:A:192:GLU:HB3	1:A:288:GLY:O	2.01	0.61
1:A:11:SER:HB3	1:A:41:PRO:HD3	1.83	0.60
1:A:192:GLU:HG3	2:A:1353:SO4:O3	2.01	0.60
1:A:13:SER:HB2	1:A:27:TRP:CH2	2.37	0.60
1:A:182:LEU:HB2	1:A:297:TYR:CE1	2.37	0.59
1:A:324:LYS:CE	2:A:1352:SO4:O1	2.51	0.59
1:A:274:PRO:HB2	1:A:277:TRP:CD1	2.38	0.59
1:A:86:GLY:O	6:A:2040:HOH:O	2.16	0.59
1:A:330:LEU:HD13	1:A:335:GLU:HB3	1.85	0.58
1:A:335:GLU:C	1:A:338:PRO:HD2	2.28	0.58
1:A:134:ASP:OD1	1:A:134:ASP:C	2.45	0.58
1:A:85:ILE:O	1:A:85:ILE:HG13	2.04	0.57
1:A:249:TYR:CE1	1:A:256:GLN:HG3	2.39	0.57
1:A:85:ILE:HG12	1:A:90:PHE:HZ	1.69	0.57
1:A:211:LYS:HD3	1:A:291:ILE:HD12	1.87	0.57
1:A:26:ALA:HB1	1:A:213:TYR:CZ	2.40	0.56
1:A:182:LEU:CD1	1:A:184:SER:N	2.68	0.56
1:A:29:GLU:OE1	1:A:29:GLU:CA	2.55	0.55
1:A:101:LEU:O	1:A:118:SER:HB2	2.06	0.55
1:A:249:TYR:CD1	1:A:256:GLN:HG3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:LEU:O	1:A:85:ILE:CG2	2.53	0.55
1:A:237:ASP:O	1:A:238:ARG:HB2	2.07	0.54
1:A:275:MET:O	1:A:276:TYR:HB2	2.07	0.54
1:A:82:GLN:HE21	1:A:126:HIS:HA	1.73	0.54
1:A:177:ARG:NH2	6:A:2066:HOH:O	2.41	0.53
1:A:211:LYS:HB3	1:A:291:ILE:HD12	1.91	0.53
1:A:31:GLN:C	1:A:285:LEU:HD12	2.33	0.53
1:A:243:ASP:OD1	1:A:243:ASP:C	2.52	0.53
1:A:77:ASP:C	1:A:77:ASP:OD1	2.53	0.52
1:A:12:GLY:HA2	1:A:262:GLU:HB3	1.92	0.52
1:A:87:ASN:O	1:A:87:ASN:ND2	2.44	0.51
1:A:143:ARG:NH1	2:A:1353:SO4:O1	2.42	0.51
1:A:104:ASP:OD1	1:A:104:ASP:C	2.53	0.51
1:A:233:HIS:CG	1:A:333:PRO:HB3	2.45	0.51
1:A:131:LYS:O	1:A:135:ILE:HD12	2.11	0.51
1:A:192:GLU:OE1	1:A:286:ASN:CA	2.58	0.51
1:A:251:ARG:C	1:A:253:PRO:HD2	2.36	0.50
1:A:87:ASN:ND2	1:A:87:ASN:C	2.69	0.49
1:A:182:LEU:HD11	1:A:184:SER:CA	2.43	0.49
1:A:67:THR:O	1:A:68:ASN:HB2	2.12	0.49
1:A:243:ASP:OD1	1:A:245:ASP:N	2.41	0.49
1:A:182:LEU:HD12	1:A:183:THR:C	2.38	0.49
1:A:17:ARG:HE	1:A:44:ARG:NE	2.10	0.48
1:A:132:LEU:HD11	1:A:191:MET:HG2	1.94	0.48
1:A:146:LEU:C	1:A:146:LEU:HD23	2.38	0.48
1:A:192:GLU:OE1	1:A:286:ASN:HA	2.14	0.48
1:A:340:LEU:O	1:A:344:ILE:HD12	2.14	0.47
1:A:154:VAL:HG23	1:A:159:VAL:CG2	2.43	0.47
1:A:183:THR:HB	1:A:296:TRP:NE1	2.29	0.47
1:A:306:ILE:HG23	1:A:307:GLU:N	2.29	0.47
1:A:112:GLN:N	1:A:112:GLN:CD	2.72	0.46
1:A:202:GLU:O	1:A:275:MET:HE3	2.14	0.46
1:A:47:GLN:HG2	6:A:2021:HOH:O	2.15	0.46
1:A:182:LEU:HD12	1:A:182:LEU:O	2.14	0.46
1:A:77:ASP:OD1	1:A:79:GLU:N	2.48	0.46
1:A:85:ILE:HG12	1:A:90:PHE:CZ	2.49	0.46
1:A:251:ARG:O	1:A:253:PRO:HD2	2.16	0.45
1:A:335:GLU:O	1:A:338:PRO:HD2	2.15	0.45
1:A:17:ARG:HH21	1:A:44:ARG:HE	1.63	0.45
1:A:160:MET:CE	1:A:163:LEU:HD12	2.28	0.45
1:A:173:GLN:O	1:A:177:ARG:CG	2.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:THR:HG22	1:A:150:LEU:N	2.31	0.44
1:A:160:MET:N	1:A:160:MET:HE3	2.32	0.44
1:A:192:GLU:OE1	1:A:287:GLY:N	2.51	0.44
1:A:13:SER:HB2	1:A:27:TRP:HH2	1.83	0.44
1:A:49:ASP:HA	1:A:50:PRO:HD2	1.88	0.44
1:A:211:LYS:O	1:A:290:THR:HA	2.17	0.44
1:A:45:LEU:HD22	1:A:49:ASP:OD2	2.18	0.43
1:A:210:ILE:HD11	1:A:293:VAL:HG23	2.00	0.43
1:A:185:ASN:HA	1:A:294:ASN:O	2.18	0.43
1:A:308:TYR:HA	1:A:309:PRO:C	2.43	0.43
1:A:82:GLN:HE21	1:A:126:HIS:CB	2.32	0.43
1:A:119:ASN:HA	6:A:2050:HOH:O	2.18	0.43
1:A:211:LYS:CB	1:A:291:ILE:HD12	2.48	0.43
1:A:47:GLN:HG3	1:A:169:TRP:NE1	2.33	0.43
1:A:222:ASP:OD1	1:A:222:ASP:C	2.62	0.43
1:A:146:LEU:HD23	1:A:146:LEU:O	2.18	0.42
1:A:207:PHE:CZ	1:A:292:THR:HG21	2.54	0.42
1:A:150:LEU:HD12	1:A:150:LEU:HA	1.37	0.42
1:A:182:LEU:HA	1:A:297:TYR:CD1	2.55	0.42
1:A:207:PHE:HB3	1:A:271:LEU:HB3	2.02	0.42
1:A:349:ASN:OD1	1:A:349:ASN:C	2.63	0.41
1:A:29:GLU:C	1:A:31:GLN:N	2.76	0.41
1:A:207:PHE:CE2	1:A:292:THR:HG21	2.56	0.41
1:A:72:PRO:HG2	1:A:164:GLY:HA3	2.02	0.41
1:A:78:LEU:HD23	1:A:78:LEU:HA	1.86	0.41
1:A:173:GLN:O	1:A:177:ARG:HG3	2.21	0.41
1:A:124:LYS:CB	1:A:127:GLU:HG3	2.46	0.41
1:A:182:LEU:HD12	1:A:183:THR:N	2.34	0.40
1:A:182:LEU:CD1	1:A:184:SER:CA	3.00	0.40
1:A:186:LEU:HD23	1:A:188:LEU:HD21	2.02	0.40
1:A:276:TYR:CD2	1:A:313:HIS:HB2	2.56	0.40
1:A:173:GLN:O	1:A:177:ARG:HG2	2.21	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:344:ILE:CD1	1:A:344:ILE:CD1[8_665]	2.02	0.18

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	294/352 (84%)	283 (96%)	10 (3%)	1 (0%)	36	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	ASP

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	A	281/306 (92%)	254 (90%)	27 (10%)	8	12

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	29	GLU
1	A	40	ARG
1	A	51	ARG
1	A	85	ILE
1	A	87	ASN
1	A	104	ASP
1	A	108	MET
1	A	135	ILE
1	A	150	LEU

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Mol	Chain	Res	Type
1	A	151	ASN
1	A	153	THR
1	A	160	MET
1	A	172	LYS
1	A	177	ARG
1	A	182	LEU
1	A	186	LEU
1	A	188	LEU
1	A	217	ILE
1	A	258	VAL
1	A	267	PRO
1	A	275	MET
1	A	302	THR
1	A	304	LYS
1	A	305	ARG
1	A	306	ILE
1	A	344	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
1	A	87	ASN
1	A	126	HIS
1	A	151	ASN
1	A	254	ASN
1	A	257	ASN
1	A	294	ASN
1	A	332	ASN
1	A	341	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 for Mogul analysis.

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	1353	-	4,4,4	0.35	0	6,6,6	0.10	0
3	GOL	A	1356	-	5,5,5	0.28	0	5,5,5	0.31	0
3	GOL	A	1357	-	5,5,5	0.56	0	5,5,5	0.28	0
4	DZA	A	1358	5	10,10,10	7.33	1 (10%)	12,12,12	3.89	7 (58%)
2	SO4	A	1351	-	4,4,4	0.38	0	6,6,6	0.09	0
2	SO4	A	1350	-	4,4,4	0.38	0	6,6,6	0.08	0
2	SO4	A	1354	-	4,4,4	0.45	0	6,6,6	0.19	0
2	SO4	A	1352	-	4,4,4	0.30	0	6,6,6	0.29	0
3	GOL	A	1355	-	5,5,5	0.49	0	5,5,5	0.24	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
4	DZA	A	1358	5	-	1/9/9/9	-
3	GOL	A	1355	-	-	0/4/4/4	-
3	GOL	A	1356	-	-	0/4/4/4	-
3	GOL	A	1357	-	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1358	DZA	N03-N04	-22.92	1.12	1.42

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1358	DZA	C06-N04-N03	-8.22	101.56	109.03
4	A	1358	DZA	C02-N03-N04	-6.04	114.32	120.29
4	A	1358	DZA	C05-N04-C06	4.38	117.01	110.98
4	A	1358	DZA	C05-N04-N03	-3.83	105.54	109.03
4	A	1358	DZA	C07-C08-C09	-3.81	103.55	113.67
4	A	1358	DZA	C08-C07-C02	-3.63	105.20	112.67
4	A	1358	DZA	O11-C09-C08	-3.46	112.12	123.09

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1358	DZA	C07-C08-C09-O10

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1353	SO4	2	0
2	A	1352	SO4	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 ⓘ

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	339/352 (96%)	0.70	41 (12%) 8 6	31, 57, 90, 98	26 (7%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	17	ARG	6.3
1	A	117	ARG	5.3
1	A	136	GLN	5.1
1	A	113	ASN	5.0
1	A	118	SER	4.9
1	A	303	PRO	4.1
1	A	115	LYS	4.1
1	A	12	GLY	4.0
1	A	135	ILE	3.8
1	A	116	PRO	3.6
1	A	286	ASN	3.6
1	A	183	THR	3.6
1	A	306	ILE	3.6
1	A	103	TYR	3.5
1	A	296	TRP	3.5
1	A	137	GLN	3.5
1	A	185	ASN	3.2
1	A	19	GLU	3.1
1	A	300	ALA	3.1
1	A	16	PRO	3.0
1	A	114	PHE	2.9
1	A	139	GLY	2.8
1	A	155	GLY	2.8
1	A	156	ARG	2.8
1	A	162	PHE	2.7
1	A	302	THR	2.7
1	A	11	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	304	LYS	2.6
1	A	51	ARG	2.5
1	A	151	ASN	2.5
1	A	152	ASP	2.4
1	A	138	ARG	2.4
1	A	129	VAL	2.4
1	A	305	ARG	2.4
1	A	92	VAL	2.3
1	A	332	ASN	2.2
1	A	100	PHE	2.1
1	A	154	VAL	2.1
1	A	142	GLU	2.1
1	A	109	ALA	2.0
1	A	13	SER	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($\text{\AA}^2$)	Q<0.9
2	SO4	A	1354	5/5	0.60	0.12	100,100,100,100	5
3	GOL	A	1355	6/6	0.68	0.33	98,100,100,100	0
2	SO4	A	1353	5/5	0.75	0.15	100,100,100,100	5
2	SO4	A	1352	5/5	0.79	0.15	100,100,100,100	5
3	GOL	A	1357	6/6	0.83	0.15	65,67,68,69	0
2	SO4	A	1350	5/5	0.86	0.16	80,81,81,82	5
3	GOL	A	1356	6/6	0.88	0.21	80,81,81,83	0
2	SO4	A	1351	5/5	0.92	0.16	91,91,91,92	5
4	DZA	A	1358	11/11	0.92	0.10	41,45,47,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	ZN	A	1359	1/1	0.99	0.03	42,42,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.