



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:09 AM UTC

PDB ID : 1USU / pdb_00001usu
Title : The Structure of the complex between Aha1 and HSP90
Authors : Meyer, P.; Roe, S.M.; Pearl, L.H.
Deposited on : 2003-12-01
Resolution : 2.15 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

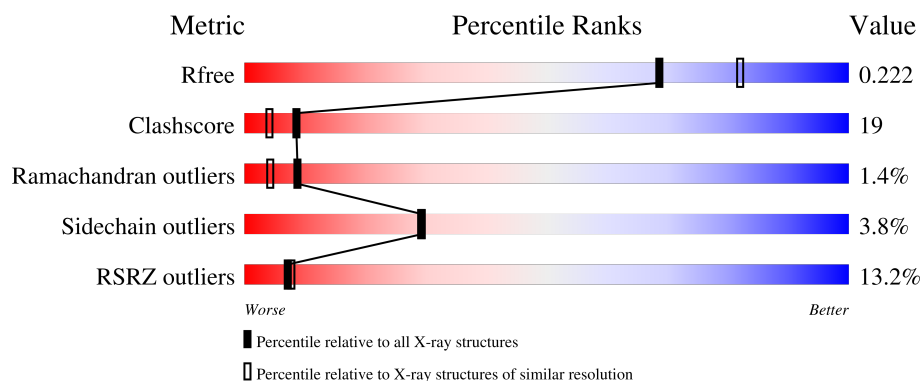
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	260	5% 67% 25% • 5%
2	B	170	21% 37% 36% • 22%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3301 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 HEAT SHOCK PROTEIN HSP82.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	0	0	0
			2031	1314	327	388	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	481	SER	ALA	conflict	UNP P02829

- Molecule 2 is a protein called AHA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	132	Total	C	N	O	S	0	0	0
			1037	666	170	199	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	28	ILE	LEU	conflict	UNP Q12449

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	199	Total	O	0	0
			199	199		
4	B	28	Total	O	0	0
			28	28		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.94Å 37.92Å 111.26Å 90.00° 98.40° 90.00°	Depositor
Resolution (Å)	31.79 – 2.15 31.79 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.1 (31.79-2.15) 99.1 (31.79-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.75 (at 2.16Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.213 , 0.268 0.216 , 0.222	Depositor DCC
R_{free} test set	1352 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.136	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3301	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.39% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	1.05	4/2074 (0.2%)	1.17	17/2802 (0.6%)
2	B	0.60	1/1052 (0.1%)	1.12	8/1410 (0.6%)
All	All	0.92	5/3126 (0.2%)	1.15	25/4212 (0.6%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	512	GLN	CA-C	6.55	1.61	1.52
1	A	473	TYR	C-O	-5.80	1.16	1.23
1	A	364	PHE	CA-C	5.54	1.60	1.52
2	B	95	VAL	CA-CB	-5.47	1.49	1.54
1	A	509	ALA	CA-CB	-5.13	1.45	1.53

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	152	ASN	N-CA-C	-8.93	102.56	113.19
1	A	485	SER	CA-C-N	8.14	127.78	119.56
1	A	485	SER	C-N-CA	8.14	127.78	119.56
2	B	28	ILE	N-CA-C	7.59	118.51	111.45
1	A	373	ASP	N-CA-C	7.17	122.32	113.50
1	A	460	TYR	N-CA-C	-7.04	103.53	111.07
1	A	355	GLU	N-CA-C	-6.93	101.59	110.53
1	A	272	PRO	N-CA-CB	6.83	110.52	103.00
2	B	129	SER	N-CA-C	6.58	118.54	111.36
2	B	95	VAL	N-CA-C	-6.58	101.63	108.15
2	B	97	GLU	N-CA-C	6.21	118.41	107.61
2	B	103	GLU	N-CA-C	-5.87	101.31	110.17
1	A	435	ASN	N-CA-C	5.76	120.43	113.17
1	A	323	ILE	CA-C-N	5.76	125.73	120.03

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	323	ILE	C-N-CA	5.76	125.73	120.03
1	A	284	ILE	N-CA-C	5.73	117.00	109.21
2	B	99	ALA	N-CA-C	5.61	117.50	109.14
1	A	299	ASP	N-CA-C	-5.59	103.11	110.43
2	B	106	SER	N-CA-C	5.58	117.44	111.36
1	A	431	GLU	N-CA-C	5.53	121.05	113.97
1	A	446	ASN	N-CA-C	-5.51	103.11	110.55
1	A	300	TRP	N-CA-C	-5.49	105.94	113.30
1	A	331	LEU	N-CA-C	5.26	117.44	111.02
1	A	471	ILE	N-CA-C	-5.12	101.03	108.36
1	A	421	PHE	N-CA-C	5.08	119.02	112.41

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	2031	0	2014	50	0
2	B	1037	0	1045	67	0
3	A	6	0	8	2	0
4	A	199	0	0	9	1
4	B	28	0	0	1	0
All	All	3301	0	3067	116	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 19.

All (116) 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)
2:B:80:ASP:HA	2:B:86:ALA:HA	1.55	0.88
2:B:14:LYS:HE3	2:B:150:HIS:HB3	1.59	0.82
2:B:134:LYS:HA	2:B:137:GLN:HE21	1.45	0.79
2:B:40:LYS:HG2	2:B:79:VAL:HG13	1.62	0.79

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:LEU:HD22	1:A:391:VAL:HG12	1.65	0.76
2:B:28:ILE:HG23	2:B:44:ILE:HD13	1.67	0.75
1:A:315:LEU:HD22	1:A:391:VAL:CG1	2.17	0.74
2:B:57:ASN:HD22	2:B:57:ASN:N	1.83	0.74
2:B:56:VAL:C	2:B:57:ASN:HD22	1.96	0.72
1:A:341:ILE:HG12	4:A:2049:HOH:O	1.91	0.70
2:B:28:ILE:CG2	2:B:44:ILE:HG21	2.22	0.70
1:A:527:ASP:C	4:A:2195:HOH:O	2.34	0.69
2:B:44:ILE:HD11	2:B:135:LEU:HD21	1.73	0.69
2:B:32:GLU:HB3	2:B:43:LYS:CB	2.23	0.68
1:A:318:ARG:NH1	4:A:2042:HOH:O	1.99	0.68
1:A:355:GLU:O	1:A:356:ASP:HB2	1.95	0.67
2:B:58:GLN:O	2:B:59:ARG:HG3	1.94	0.67
1:A:340:ASN:OD1	1:A:366:LYS:HE2	1.95	0.66
1:A:394:LYS:HE3	4:B:2012:HOH:O	1.97	0.65
2:B:80:ASP:CA	2:B:86:ALA:HA	2.25	0.65
1:A:347:ARG:NH1	1:A:372:GLU:OE2	2.31	0.64
1:A:385:GLN:O	1:A:385:GLN:HG2	1.96	0.64
2:B:147:LEU:C	2:B:149:THR:H	2.06	0.64
2:B:57:ASN:HB3	2:B:59:ARG:NH2	2.13	0.63
1:A:479:LEU:C	1:A:479:LEU:HD13	2.23	0.63
3:A:1528:GOL:H31	4:A:2127:HOH:O	1.97	0.63
1:A:475:THR:HG21	1:A:507:GLU:HG3	1.81	0.62
2:B:79:VAL:O	2:B:80:ASP:HB3	1.98	0.62
2:B:123:ALA:O	2:B:126:LEU:HB3	2.00	0.62
2:B:29:VAL:HG13	2:B:45:LYS:O	2.00	0.61
2:B:28:ILE:HG22	2:B:44:ILE:HG21	1.80	0.61
2:B:116:GLU:HA	2:B:120:LEU:HD12	1.82	0.60
2:B:63:VAL:HG11	2:B:151:GLY:HA2	1.83	0.60
2:B:61:GLY:O	2:B:62:LYS:HG3	2.02	0.60
2:B:58:GLN:OE1	2:B:63:VAL:HG13	2.03	0.59
2:B:32:GLU:HB3	2:B:43:LYS:HB3	1.86	0.58
2:B:61:GLY:C	2:B:62:LYS:HG3	2.27	0.58
2:B:76:GLU:HG2	2:B:90:GLU:HG2	1.85	0.57
2:B:91:GLY:HA2	2:B:114:PHE:CD1	2.39	0.57
2:B:149:THR:HG22	2:B:150:HIS:CE1	2.39	0.57
2:B:59:ARG:HG3	2:B:59:ARG:HH21	1.70	0.57
1:A:339:ASN:HA	1:A:352:ASP:O	2.05	0.57
1:A:300:TRP:NE1	4:A:2027:HOH:O	2.21	0.56
1:A:402:GLU:O	1:A:406:GLU:HG2	2.06	0.56
2:B:32:GLU:HB3	2:B:43:LYS:HB2	1.88	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:479:LEU:HD11	1:A:483:GLU:OE2	2.06	0.56
1:A:356:ASP:O	1:A:393:ARG:HD3	2.07	0.55
1:A:394:LYS:O	1:A:398:LYS:HG3	2.07	0.54
2:B:57:ASN:N	2:B:57:ASN:ND2	2.49	0.54
1:A:374:LEU:HD12	1:A:375:PRO:HD2	1.89	0.54
2:B:50:ILE:HG12	2:B:71:ILE:HG12	1.88	0.54
2:B:41:TYR:C	2:B:41:TYR:CD2	2.86	0.54
1:A:275:PRO:HG2	1:A:278:THR:CG2	2.37	0.54
2:B:151:GLY:C	2:B:153:ASP:H	2.16	0.54
2:B:80:ASP:HA	2:B:86:ALA:CA	2.35	0.53
2:B:56:VAL:CG1	2:B:63:VAL:HG12	2.39	0.53
1:A:275:PRO:HG2	1:A:278:THR:HG23	1.90	0.53
1:A:445:TYR:OH	1:A:506:ASP:OD1	2.20	0.53
1:A:309:PHE:CZ	1:A:399:LYS:HG3	2.44	0.53
1:A:399:LYS:NZ	4:A:2077:HOH:O	2.31	0.52
1:A:526:LYS:O	1:A:527:ASP:O	2.28	0.51
2:B:39:LYS:HB2	2:B:39:LYS:NZ	2.26	0.51
1:A:318:ARG:CZ	4:A:2042:HOH:O	2.51	0.51
1:A:318:ARG:NH2	4:A:2042:HOH:O	2.40	0.51
1:A:345:VAL:HG11	1:A:376:LEU:HD21	1.93	0.51
2:B:59:ARG:HG3	2:B:59:ARG:NH2	2.25	0.50
1:A:346:ARG:NH1	1:A:372:GLU:O	2.44	0.50
2:B:114:PHE:O	2:B:115:LYS:C	2.54	0.50
1:A:320:ILE:O	1:A:320:ILE:HG23	2.11	0.50
1:A:355:GLU:O	1:A:356:ASP:CB	2.60	0.50
1:A:440:ALA:HB2	1:A:513:LEU:HD21	1.93	0.50
2:B:45:LYS:HD2	2:B:76:GLU:CD	2.37	0.50
2:B:80:ASP:CB	2:B:86:ALA:HA	2.41	0.50
2:B:151:GLY:C	2:B:153:ASP:N	2.70	0.50
1:A:476:GLY:HA3	1:A:482:VAL:HG13	1.94	0.49
1:A:333:GLU:O	1:A:334:SER:C	2.56	0.49
2:B:41:TYR:HA	2:B:126:LEU:HD21	1.95	0.49
2:B:87:LEU:HD12	2:B:88:PRO:HD2	1.94	0.49
1:A:458:THR:HB	3:A:1528:GOL:H2	1.94	0.48
2:B:17:ILE:HG13	2:B:21:LYS:HD2	1.95	0.48
2:B:23:TYR:CE2	2:B:27:LYS:HG3	2.48	0.48
1:A:514:LYS:NZ	2:B:110:ASP:OD1	2.45	0.47
2:B:41:TYR:CA	2:B:126:LEU:HD21	2.44	0.47
2:B:58:GLN:C	2:B:59:ARG:HG3	2.39	0.47
1:A:476:GLY:HA3	1:A:482:VAL:CG1	2.44	0.47
1:A:315:LEU:CD2	1:A:391:VAL:HG12	2.39	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:77:GLY:O	2:B:88:PRO:HA	2.15	0.47
2:B:91:GLY:HA3	2:B:112:SER:O	2.15	0.46
2:B:63:VAL:HG12	2:B:63:VAL:O	2.14	0.46
1:A:340:ASN:HB3	1:A:366:LYS:HD3	1.98	0.46
2:B:17:ILE:O	2:B:21:LYS:HD2	2.16	0.46
2:B:87:LEU:HD12	2:B:88:PRO:CD	2.45	0.46
1:A:314:GLN:HE21	2:B:64:ILE:HB	1.81	0.45
2:B:87:LEU:HA	2:B:88:PRO:HD3	1.78	0.45
1:A:409:GLU:HG3	4:A:2086:HOH:O	2.16	0.45
2:B:40:LYS:HE2	2:B:79:VAL:CG1	2.47	0.45
1:A:339:ASN:CB	1:A:352:ASP:O	2.65	0.44
1:A:460:TYR:C	1:A:460:TYR:CD2	2.96	0.44
2:B:23:TYR:C	2:B:23:TYR:CD2	2.95	0.44
2:B:25:LYS:HB3	2:B:25:LYS:HE2	1.75	0.44
1:A:305:TYR:OH	1:A:406:GLU:HG3	2.19	0.43
2:B:28:ILE:HG21	2:B:73:VAL:HG11	2.00	0.43
1:A:477:GLU:O	1:A:477:GLU:HG2	2.18	0.43
2:B:48:SER:O	2:B:49:SER:HB3	2.18	0.43
2:B:79:VAL:O	2:B:80:ASP:CB	2.65	0.43
2:B:122:GLU:C	2:B:125:PRO:HD2	2.43	0.43
2:B:12:VAL:N	2:B:58:GLN:HG3	2.35	0.42
2:B:40:LYS:HG2	2:B:79:VAL:CG1	2.43	0.42
1:A:315:LEU:HD22	1:A:391:VAL:HG11	1.97	0.42
1:A:377:ASN:O	1:A:378:LEU:HG	2.19	0.42
2:B:28:ILE:O	2:B:31:VAL:HG23	2.19	0.41
2:B:32:GLU:HA	2:B:43:LYS:HA	2.02	0.41
1:A:339:ASN:CA	1:A:352:ASP:O	2.68	0.41
1:A:309:PHE:CE1	1:A:399:LYS:HG3	2.57	0.40
1:A:314:GLN:H	1:A:314:GLN:HG2	1.45	0.40
2:B:85:SER:O	2:B:86:ALA:HB2	2.22	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)
4:A:2171:HOH:O	4:A:2196:HOH:O[2_646]	2.09	0.11

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	240/260 (92%)	232 (97%)	7 (3%)	1 (0%)	30	26
2	B	124/170 (73%)	107 (86%)	13 (10%)	4 (3%)	3	0
All	All	364/430 (85%)	339 (93%)	20 (6%)	5 (1%)	9	4

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	121	SER
2	B	148	ALA
1	A	333	GLU
2	B	64	ILE
2	B	79	VAL

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	226/241 (94%)	219 (97%)	7 (3%)	35	37
2	B	115/149 (77%)	109 (95%)	6 (5%)	21	17
All	All	341/390 (87%)	328 (96%)	13 (4%)	29	29

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	306	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	314	GLN
1	A	333	GLU
1	A	353	GLU
1	A	453	GLU
1	A	469	LYS
1	A	507	GLU
2	B	26	GLN
2	B	39	LYS
2	B	41	TYR
2	B	57	ASN
2	B	112	SER
2	B	119	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	314	GLN
1	A	395	ASN
1	A	405	ASN
1	A	413	GLN
2	B	26	GLN
2	B	57	ASN
2	B	137	GLN
2	B	152	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 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
3	GOL	A	1528	-	5,5,5	0.83	0	5,5,5	0.98	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
3	GOL	A	1528	-	-	0/4/4/4	-

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
3	A	1528	GOL	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

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	246/260 (94%)	0.20	14 (5%) 29 33	24, 34, 60, 90	0
2	B	132/170 (77%)	1.45	36 (27%) 1 1	37, 69, 95, 100	0
All	All	378/430 (87%)	0.63	50 (13%) 7 8	24, 42, 87, 100	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	329	PHE	5.5
1	A	272	PRO	4.5
1	A	375	PRO	4.5
2	B	148	ALA	4.4
1	A	527	ASP	4.4
2	B	149	THR	4.0
2	B	30	GLY	3.9
2	B	79	VAL	3.8
2	B	77	GLY	3.2
2	B	88	PRO	3.2
2	B	120	LEU	3.1
1	A	355	GLU	3.1
2	B	87	LEU	3.0
2	B	16	CYS	3.0
2	B	29	VAL	2.9
2	B	153	ASP	2.9
1	A	354	ALA	2.9
1	A	334	SER	2.9
2	B	85	SER	2.9
2	B	31	VAL	2.8
2	B	17	ILE	2.8
1	A	356	ASP	2.8
2	B	61	GLY	2.8
2	B	78	HIS	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	332	PHE	2.6
2	B	19	TRP	2.6
2	B	46	SER	2.5
2	B	89	PHE	2.5
2	B	13	ASP	2.5
2	B	110	ASP	2.5
1	A	295	SER	2.5
2	B	40	LYS	2.5
2	B	117	THR	2.4
2	B	86	ALA	2.4
2	B	76	GLU	2.4
2	B	64	ILE	2.3
2	B	12	VAL	2.3
2	B	118	SER	2.3
1	A	378	LEU	2.3
2	B	41	TYR	2.3
1	A	273	THR	2.2
1	A	376	LEU	2.2
2	B	132	LEU	2.2
2	B	150	HIS	2.2
2	B	75	ILE	2.2
2	B	22	GLU	2.2
2	B	126	LEU	2.2
2	B	114	PHE	2.1
2	B	20	ALA	2.1
1	A	340	ASN	2.1

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
3	GOL	A	1528	6/6	0.87	0.19	46,48,49,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.