Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:09 PM GMT

PDB ID	:	1SBS
Title	:	CRYSTAL STRUCTURE OF AN ANTI-HCG FAB
Authors	:	Fotinou, C.; Beauchamp, J.; Emsley, P.; Dehaan, A.; Schielen, W.J.G.; Bos, E.; Isaacs, N.W.
Deposited on	:	1998-04-08
Resolution	:	2.00 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see references ⓘ) were used in the production of this report:

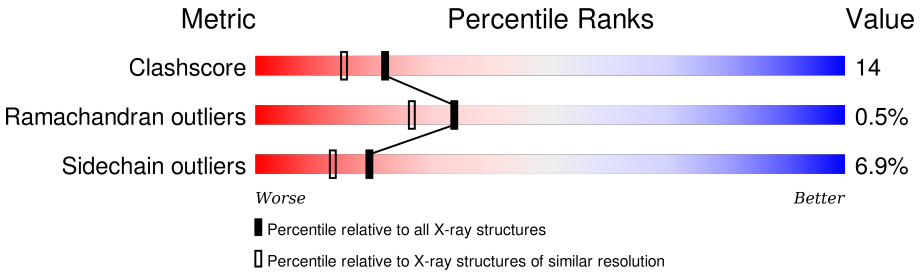
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	H	222	80% 16%
2	L	220	74% 23%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4360 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 MONOCLONAL ANTIBODY 3A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	222	Total	C	N	O	S	0	0	0
			1696	1069	281	337	9			

There are 23 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	3	ASN	LYS	CONFLICT	GB 1042226
H	20	LEU	VAL	CONFLICT	GB 1042226
H	28	THR	ALA	CONFLICT	GB 1042226
H	31	ASN	TYR	CONFLICT	GB 1042226
H	43	LYS	ARG	CONFLICT	GB 1042226
H	48	VAL	ILE	CONFLICT	GB 1042226
H	50	ASP	GLU	CONFLICT	GB 1042226
H	53	LEU	PHE	CONFLICT	GB 1042226
H	59	ALA	-	INSERTION	GB 1042226
H	61	LEU	HIS	CONFLICT	GB 1042226
H	101	GLY	GLU	CONFLICT	GB 1042226
H	102	ALA	GLY	CONFLICT	GB 1042226
H	103	TYR	ILE	CONFLICT	GB 1042226
H	105	ARG	-	INSERTION	GB 1042226
H	106	TYR	-	INSERTION	GB 1042226
H	107	ASP	-	INSERTION	GB 1042226
H	109	ALA	PRO	CONFLICT	GB 1042226
H	110	MET	PHE	CONFLICT	GB 1042226
H	111	ASP	ALA	CONFLICT	GB 1042226
H	118	SER	LEU	CONFLICT	GB 1042226
H	123	SER	ALA	CONFLICT	GB 1042226
H	197	PRO	THR	CONFLICT	GB 1042226
H	198	ARG	TRP	CONFLICT	GB 1042226

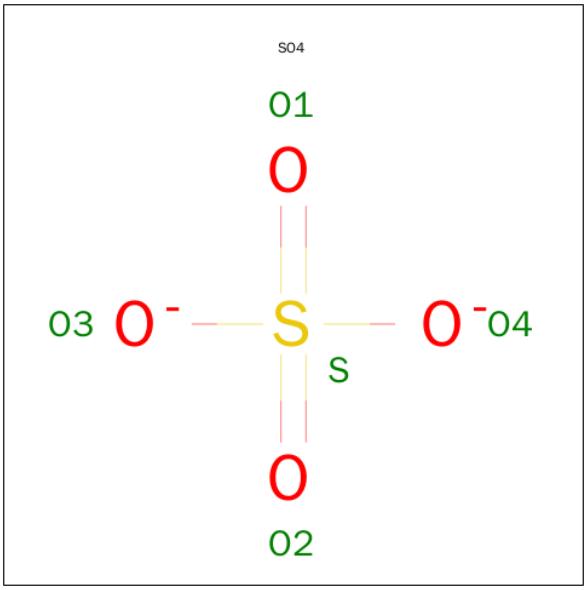
- Molecule 2 is a protein called MONOCLONAL ANTIBODY 3A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	220	Total	C	N	O	S	0	0	0
			1715	1071	282	353	9			

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	5	SER	THR	CONFLICT	GB 1407830
L	15	VAL	ALA	CONFLICT	GB 1407830
L	22	THR	SER	CONFLICT	GB 1407830
L	31	TYR	ASN	CONFLICT	GB 1407830
L	32	SER	THR	CONFLICT	GB 1407830
L	33	SER	TRP	CONFLICT	GB 1407830
L	34	ASN	THR	CONFLICT	GB 1407830
L	35	GLN	ARG	CONFLICT	GB 1407830
L	36	MET	LYS	CONFLICT	GB 1407830
L	85	GLU	GLN	CONFLICT	GB 1407830
L	95	GLN	LYS	CONFLICT	GB 1407830
L	97	TYR	-	INSERTION	GB 1407830
L	98	HIS	-	INSERTION	GB 1407830
L	101	PRO	ASN	CONFLICT	GB 1407830
L	?	-	PRO	DELETION	GB 1407830
L	106	SER	GLY	CONFLICT	GB 1407830
L	109	LYS	GLN	CONFLICT	GB 1407830

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	H	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	H	452	Total O 452 452	0	0
4	L	492	Total O 492 492	0	0

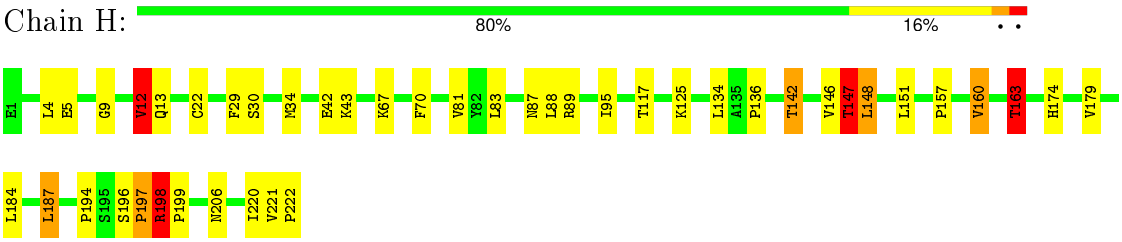
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

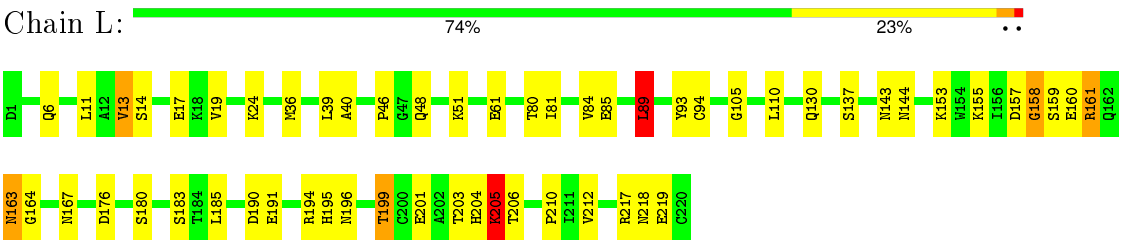
- Molecule 1: MONOCLONAL ANTIBODY 3A2

Chain H:



- Molecule 2: MONOCLONAL ANTIBODY 3A2

Chain L:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	74.80Å 74.80Å 198.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.00	Depositor
% Data completeness (in resolution range)	99.0 (20.00-2.00)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.180 , 0.240	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4360	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

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	H	0.55	1/1740 (0.1%)	1.21	11/2374 (0.5%)
2	L	0.53	0/1755	1.22	11/2382 (0.5%)
All	All	0.54	1/3495 (0.0%)	1.22	22/4756 (0.5%)

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
2	L	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	222	PRO	N-CD	6.56	1.57	1.47

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	L	164	GLY	CA-C-O	9.39	137.51	120.60
2	L	163	ASN	CA-C-N	8.95	134.11	116.20
1	H	198	ARG	CB-CG-CD	8.83	134.55	111.60
2	L	164	GLY	O-C-N	-7.44	110.79	122.70
1	H	89	ARG	NE-CZ-NH1	6.57	123.59	120.30
1	H	198	ARG	CG-CD-NE	6.47	125.38	111.80
1	H	12	VAL	CB-CA-C	-6.36	99.31	111.40
2	L	89	LEU	CA-CB-CG	6.27	129.72	115.30
2	L	176	ASP	CB-CG-OD1	5.97	123.68	118.30
2	L	157	ASP	CA-C-N	-5.88	104.43	116.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	147	THR	N-CA-CB	5.81	121.34	110.30
2	L	161	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	H	198	ARG	NE-CZ-NH1	-5.39	117.61	120.30
1	H	163	THR	N-CA-CB	-5.39	100.07	110.30
1	H	70	PHE	CB-CG-CD1	5.31	124.52	120.80
2	L	163	ASN	N-CA-C	5.25	125.16	111.00
2	L	163	ASN	CA-C-O	-5.23	109.12	120.10
2	L	14	SER	CB-CA-C	5.16	119.91	110.10
2	L	157	ASP	CA-C-O	5.13	130.87	120.10
1	H	187	LEU	CA-CB-CG	5.12	127.08	115.30
1	H	221	VAL	CA-C-O	5.07	130.74	120.10
1	H	160	VAL	N-CA-CB	-5.05	100.39	111.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	L	163	ASN	Mainchain

5.2 Too-close contacts (i)

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	H	1696	0	1640	53	0
2	L	1715	0	1635	42	6
3	H	5	0	0	0	0
4	H	452	0	0	21	7
4	L	492	0	0	17	6
All	All	4360	0	3275	93	14

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 14.

All (93) 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:H:179:VAL:HB	4:H:668:HOH:O	1.28	1.32
1:H:146:VAL:HB	4:H:670:HOH:O	1.47	1.11
1:H:198:ARG:HG2	1:H:199:PRO:HA	1.46	0.97
1:H:198:ARG:HB3	1:H:198:ARG:HH11	1.37	0.89
1:H:194:PRO:O	1:H:197:PRO:HD2	1.72	0.88
2:L:155:LYS:HD3	2:L:159:SER:HA	1.55	0.88
2:L:36:MET:SD	4:L:483:HOH:O	2.33	0.86
2:L:36:MET:CE	4:L:483:HOH:O	2.27	0.83
1:H:194:PRO:HD2	1:H:197:PRO:HG2	1.61	0.81
1:H:9:GLY:H	1:H:117:THR:HG21	1.45	0.81
2:L:159:SER:HB2	4:L:311:HOH:O	1.84	0.78
2:L:196:ASN:HD21	2:L:218:ASN:H	1.32	0.77
2:L:130:GLN:HE22	2:L:137:SER:H	1.34	0.76
1:H:198:ARG:CB	1:H:198:ARG:HH11	1.98	0.75
1:H:197:PRO:HB3	4:H:656:HOH:O	1.88	0.74
1:H:81:VAL:HB	4:H:655:HOH:O	1.87	0.74
1:H:174:HIS:HE1	2:L:144:ASN:HD21	1.39	0.71
2:L:130:GLN:NE2	2:L:137:SER:H	1.89	0.70
2:L:201:GLU:HG2	2:L:212:VAL:HG22	1.72	0.70
1:H:117:THR:HG23	4:H:402:HOH:O	1.93	0.68
1:H:163:THR:HG21	4:H:644:HOH:O	1.97	0.65
1:H:163:THR:HG22	1:H:206:ASN:HB2	1.79	0.65
1:H:146:VAL:CB	4:H:670:HOH:O	2.23	0.64
1:H:9:GLY:H	1:H:117:THR:CG2	2.12	0.61
1:H:95:ILE:HD11	4:H:637:HOH:O	2.00	0.61
1:H:29:PHE:HZ	1:H:81:VAL:HG13	1.65	0.61
1:H:142:THR:HG23	4:H:328:HOH:O	2.00	0.60
1:H:198:ARG:CG	1:H:198:ARG:HH11	2.13	0.60
1:H:95:ILE:HD13	4:H:520:HOH:O	2.00	0.60
1:H:12:VAL:HG21	1:H:88:LEU:HD13	1.83	0.60
2:L:155:LYS:CD	2:L:159:SER:HA	2.31	0.60
2:L:48:GLN:HG3	4:L:712:HOH:O	2.02	0.59
2:L:196:ASN:ND2	2:L:218:ASN:H	1.98	0.58
2:L:80:THR:HG23	4:L:246:HOH:O	2.02	0.58
1:H:43:LYS:HD3	4:H:351:HOH:O	2.04	0.57
2:L:48:GLN:HG3	4:L:659:HOH:O	2.06	0.56
1:H:29:PHE:CZ	1:H:81:VAL:HG13	2.41	0.55
1:H:5:GLU:HB2	4:H:611:HOH:O	2.06	0.55
1:H:148:LEU:HG	1:H:220:ILE:HG21	1.88	0.55
2:L:24:LYS:HE2	4:L:489:HOH:O	2.06	0.55
1:H:34:MET:HG3	1:H:81:VAL:HG11	1.89	0.55
2:L:199:THR:HG23	4:L:658:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:174:HIS:HD2	4:H:410:HOH:O	1.89	0.54
1:H:174:HIS:CE1	2:L:144:ASN:HD21	2.24	0.54
1:H:196:SER:OG	1:H:197:PRO:HD3	2.07	0.54
1:H:194:PRO:HB2	1:H:197:PRO:CD	2.38	0.54
2:L:6:GLN:HE22	2:L:93:TYR:HA	1.74	0.53
2:L:203:THR:HG23	2:L:210:PRO:HB3	1.91	0.52
1:H:148:LEU:HD21	1:H:198:ARG:HD3	1.91	0.52
2:L:61:GLU:HG2	4:L:356:HOH:O	2.10	0.52
1:H:34:MET:CB	1:H:81:VAL:HG11	2.40	0.51
1:H:9:GLY:HA3	1:H:117:THR:HG22	1.92	0.51
1:H:198:ARG:HB3	1:H:198:ARG:NH1	2.18	0.51
1:H:87:ASN:HB3	4:H:578:HOH:O	2.11	0.51
2:L:46:PRO:HG2	2:L:89:LEU:HD13	1.93	0.50
2:L:159:SER:HB2	4:L:239:HOH:O	2.11	0.50
2:L:19:VAL:HG22	2:L:81:ILE:HB	1.94	0.50
2:L:204:HIS:O	2:L:205:LYS:C	2.50	0.49
1:H:198:ARG:NH1	4:H:653:HOH:O	2.33	0.49
2:L:143:ASN:HD22	2:L:180:SER:HB3	1.78	0.49
2:L:13:VAL:HG22	2:L:17:GLU:OE1	2.12	0.48
2:L:39:LEU:HG	2:L:40:ALA:N	2.28	0.48
2:L:160:GLU:HG3	4:L:225:HOH:O	2.14	0.47
2:L:219:GLU:HG3	4:L:411:HOH:O	2.15	0.47
1:H:194:PRO:HB2	1:H:197:PRO:HD2	1.97	0.47
2:L:191:GLU:HG3	2:L:194:ARG:NH2	2.31	0.45
1:H:67:LYS:HD2	4:H:417:HOH:O	2.16	0.45
2:L:158:GLY:HA2	4:L:654:HOH:O	2.15	0.45
1:H:13:GLN:HB2	4:H:432:HOH:O	2.16	0.45
2:L:203:THR:CG2	2:L:210:PRO:HB3	2.46	0.45
1:H:22:CYS:HB3	1:H:81:VAL:HG23	1.97	0.45
1:H:34:MET:HG3	1:H:81:VAL:CG1	2.47	0.45
1:H:136:PRO:HD3	1:H:148:LEU:HD12	1.98	0.44
1:H:197:PRO:HD3	4:H:521:HOH:O	2.18	0.44
2:L:51:LYS:HD2	4:L:673:HOH:O	2.17	0.44
2:L:130:GLN:HE22	2:L:137:SER:N	2.10	0.43
2:L:195:HIS:O	2:L:217:ARG:HD3	2.18	0.43
1:H:142:THR:HG21	4:H:280:HOH:O	2.18	0.43
2:L:190:ASP:O	2:L:194:ARG:HG3	2.19	0.43
2:L:167:ASN:HD22	2:L:183:SER:HA	1.83	0.42
2:L:159:SER:CB	4:L:239:HOH:O	2.66	0.42
1:H:146:VAL:CG1	4:H:670:HOH:O	2.61	0.42
2:L:6:GLN:NE2	2:L:94:CYS:H	2.17	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:194:PRO:C	1:H:197:PRO:HD2	2.38	0.42
1:H:147:THR:CG2	4:L:364:HOH:O	2.68	0.42
2:L:153:LYS:HE2	2:L:155:LYS:HE3	2.02	0.42
2:L:17:GLU:O	2:L:84:VAL:HG23	2.19	0.41
1:H:198:ARG:NH2	4:H:653:HOH:O	2.41	0.41
2:L:6:GLN:HE21	2:L:105:GLY:HA3	1.86	0.41
1:H:87:ASN:HD22	1:H:87:ASN:N	2.18	0.41
1:H:34:MET:CG	1:H:81:VAL:HG11	2.52	0.40
1:H:147:THR:HG22	4:L:364:HOH:O	2.20	0.40
1:H:9:GLY:N	1:H:117:THR:HG21	2.26	0.40

All (14) 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:H:654:HOH:O	4:L:453:HOH:O[6_666]	1.76	0.44
4:H:527:HOH:O	4:L:357:HOH:O[6_666]	1.93	0.27
2:L:85:GLU:OE2	2:L:158:GLY:CA[5_666]	1.97	0.23
4:H:639:HOH:O	4:L:320:HOH:O[6_666]	1.97	0.23
2:L:85:GLU:OE1	2:L:158:GLY:CA[5_666]	2.01	0.19
2:L:85:GLU:CD	2:L:158:GLY:CA[5_666]	2.08	0.12
4:H:544:HOH:O	4:H:544:HOH:O[5_676]	2.11	0.09
4:H:597:HOH:O	4:H:597:HOH:O[5_676]	2.13	0.07
4:H:326:HOH:O	4:L:650:HOH:O[6_666]	2.13	0.07
2:L:85:GLU:OE1	2:L:158:GLY:O[5_666]	2.13	0.07
2:L:85:GLU:OE2	2:L:158:GLY:N[5_666]	2.14	0.06
4:H:518:HOH:O	4:L:408:HOH:O[6_666]	2.15	0.05
4:L:404:HOH:O	4:L:469:HOH:O[5_666]	2.17	0.03
2:L:85:GLU:OE1	2:L:158:GLY:C[5_666]	2.19	0.01

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	H	220/222 (99%)	217 (99%)	3 (1%)	0	100 100
2	L	218/220 (99%)	210 (96%)	6 (3%)	2 (1%)	21 13
All	All	438/442 (99%)	427 (98%)	9 (2%)	2 (0%)	34 26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	158	GLY
2	L	205	LYS

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	H	191/191 (100%)	173 (91%)	18 (9%)	11 6
2	L	198/198 (100%)	189 (96%)	9 (4%)	34 29
All	All	389/389 (100%)	362 (93%)	27 (7%)	19 13

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

Mol	Chain	Res	Type
1	H	4	LEU
1	H	12	VAL
1	H	30	SER
1	H	42	GLU
1	H	83	LEU
1	H	125	LYS
1	H	134	LEU
1	H	142	THR
1	H	147	THR
1	H	148	LEU
1	H	151	LEU
1	H	157	PRO
1	H	160	VAL
1	H	163	THR

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Mol	Chain	Res	Type
1	H	184	LEU
1	H	187	LEU
1	H	197	PRO
1	H	198	ARG
2	L	11	LEU
2	L	13	VAL
2	L	89	LEU
2	L	110	LEU
2	L	161	ARG
2	L	185	LEU
2	L	199	THR
2	L	205	LYS
2	L	206	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	31	ASN
1	H	87	ASN
1	H	174	HIS
2	L	6	GLN
2	L	48	GLN
2	L	130	GLN
2	L	143	ASN
2	L	144	ASN
2	L	167	ASN
2	L	195	HIS
2	L	196	ASN
2	L	204	HIS

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 carbohydrates 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	SO4	H	-1	-	4,4,4	0.83	0	6,6,6	0.22	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	SO4	H	-1	-	-	0/0/0/0	0/0/0/0

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

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.