



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:06 AM GMT

PDB ID : 1T6B
Title : Crystal structure of B. anthracis Protective Antigen complexed with human Anthrax toxin receptor
Authors : Santelli, E.; Bankston, L.A.; Leppla, S.H.; Liddington, R.C.
Deposited on : 2004-05-05
Resolution : 2.50 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

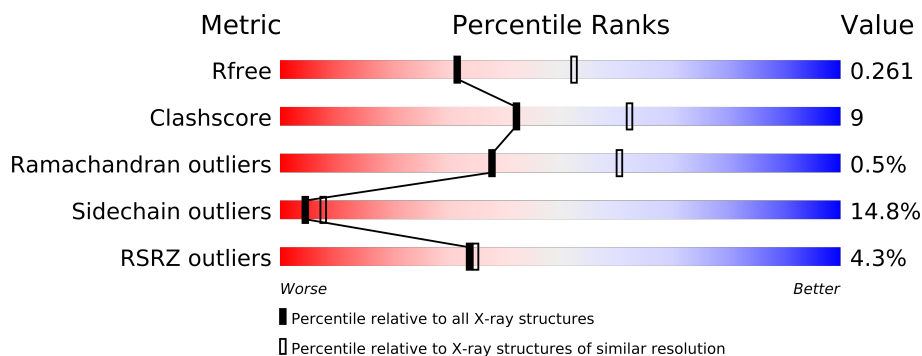
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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))
R_{free}	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	X	735	
2	Y	189	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	CA	X	801	-	X
4	MN	Y	303	-	X
5	NA	X	802	-	X
5	NA	X	803	-	X
6	PG4	X	804	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6852 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Protective Antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	676	Total	C	N	O	S	0	0	0
			5379	3378	914	1079	8			

- Molecule 2 is a protein called Anthrax toxin receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Y	170	Total	C	N	O	S	0	0	0
			1316	846	220	248	2			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Y	30	GLY	-	CLONING ARTIFACT	UNP P58335
Y	31	SER	-	CLONING ARTIFACT	UNP P58335
Y	32	HIS	-	CLONING ARTIFACT	UNP P58335
Y	33	MET	-	CLONING ARTIFACT	UNP P58335
Y	34	LEU	-	CLONING ARTIFACT	UNP P58335
Y	36	ASP	GLN	CLONING ARTIFACT	UNP P58335
Y	38	ARG	SER	CLONING ARTIFACT	UNP P58335
Y	39	GLY	CYS	CLONING ARTIFACT	UNP P58335
Y	213	SER	-	CLONING ARTIFACT	UNP P58335
Y	214	ASN	-	CLONING ARTIFACT	UNP P58335
Y	215	GLY	-	CLONING ARTIFACT	UNP P58335
Y	216	ILE	-	CLONING ARTIFACT	UNP P58335
Y	217	ALA	-	CLONING ARTIFACT	UNP P58335
Y	218	ALA	-	CLONING ARTIFACT	UNP P58335

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	X	2	Total	Ca	0	0
			2	2		

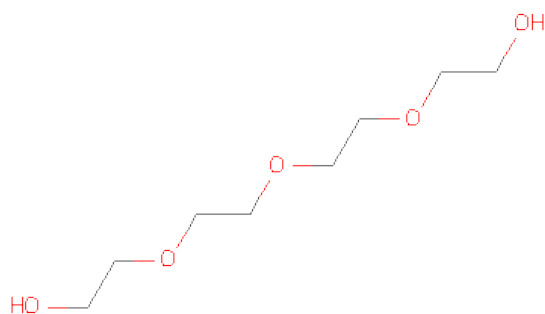
- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	Y	1	Total	Mn	0	0
			1	1		

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	X	2	Total	Na	0	0
			2	2		

- Molecule 6 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	X	1	Total	C	O	0	0
			13	8	5		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	X	119	Total	O	0	0
			119	119		
7	Y	20	Total	O	0	0
			20	20		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.16Å 94.12Å 135.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.50 29.56 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.50) 99.7 (29.56-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.207 , 0.266 0.202 , 0.261	Depositor DCC
R_{free} test set	1990 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.441	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 14.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 39610 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6852	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: NA, CA, MN, PG4

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	X	0.80	2/5474 (0.0%)	0.95	23/7418 (0.3%)
2	Y	0.79	0/1337	0.83	1/1801 (0.1%)
All	All	0.80	2/6811 (0.0%)	0.93	24/9219 (0.3%)

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	Y	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	437	ASN	CB-CG	5.86	1.64	1.51
1	X	525	GLU	CG-CD	5.73	1.60	1.51

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	512	ASP	CB-CG-OD2	8.75	126.17	118.30
1	X	595	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	X	520	ASP	CB-CG-OD2	7.37	124.93	118.30
2	Y	98	ASP	CB-CG-OD2	7.29	124.86	118.30
1	X	231	ASP	CB-CG-OD2	7.18	124.77	118.30
1	X	608	ASP	CB-CG-OD2	6.97	124.57	118.30
1	X	131	ASP	CB-CG-OD2	6.65	124.28	118.30
1	X	425	ASP	CB-CG-OD2	6.53	124.18	118.30
1	X	265	ASP	CB-CG-OD2	6.39	124.06	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	555	ASP	CB-CG-OD2	6.36	124.02	118.30
1	X	472	ASP	CB-CG-OD2	6.35	124.01	118.30
1	X	84	ASP	CB-CG-OD2	6.34	124.00	118.30
1	X	658	ASP	CB-CG-OD2	6.28	123.95	118.30
1	X	177	ASP	CB-CG-OD2	6.09	123.78	118.30
1	X	595	ARG	NE-CZ-NH2	-5.89	117.35	120.30
1	X	677	ASP	CB-CG-OD2	5.84	123.56	118.30
1	X	451	ASP	CB-CG-OD2	5.60	123.34	118.30
1	X	365	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	X	497	ASP	CB-CG-OD2	5.33	123.10	118.30
1	X	203	LEU	CB-CG-CD2	-5.27	102.05	111.00
1	X	551	ASP	CB-CG-OD2	5.26	123.03	118.30
1	X	546	ASP	CB-CG-OD2	5.21	122.98	118.30
1	X	661	ASP	CB-CG-OD2	5.18	122.97	118.30
1	X	181	ASP	CB-CG-OD1	5.03	122.82	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	Y	42	ALA	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	X	5379	0	5313	106	0
2	Y	1316	0	1348	21	0
3	X	2	0	0	0	0
4	Y	1	0	0	0	0
5	X	2	0	0	0	0
6	X	13	0	18	1	0
7	X	119	0	0	20	0
7	Y	20	0	0	2	0
All	All	6852	0	6679	120	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (120) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:X:409:ASN:HD22	1:X:410:TYR:H	1.12	0.97
1:X:327:SER:HA	7:X:915:HOH:O	1.61	0.96
1:X:521:MET:HE1	1:X:525:GLU:HB3	1.45	0.94
1:X:521:MET:CE	1:X:525:GLU:HB3	1.99	0.93
1:X:425:ASP:HA	7:X:869:HOH:O	1.69	0.90
1:X:123:GLU:O	1:X:125:PRO:HD3	1.75	0.84
1:X:345:THR:HG22	1:X:443:GLU:OE2	1.79	0.82
1:X:496:LYS:HB3	7:X:913:HOH:O	1.82	0.80
1:X:654:GLU:OE1	2:Y:51:LYS:NZ	2.21	0.73
1:X:236:PHE:O	1:X:240:THR:HB	1.90	0.72
1:X:629:LEU:HD12	1:X:631:ILE:HD11	1.71	0.72
1:X:656:ILE:HD12	2:Y:115:VAL:HG11	1.72	0.72
1:X:175:VAL:HA	7:X:839:HOH:O	1.93	0.68
1:X:338:LEU:O	1:X:338:LEU:HD23	1.93	0.68
1:X:240:THR:HG21	1:X:242:ARG:HH11	1.59	0.68
1:X:152:GLN:HB3	7:X:827:HOH:O	1.95	0.67
1:X:521:MET:HE2	1:X:522:THR:O	1.95	0.66
1:X:298:THR:HG22	7:X:902:HOH:O	1.96	0.66
1:X:263:HIS:HD2	7:X:877:HOH:O	1.79	0.63
1:X:46:PRO:HD2	1:X:49:GLU:OE2	2.01	0.61
1:X:593:ASP:OD1	1:X:595:ARG:HG2	2.02	0.60
1:X:231:ASP:HB2	1:X:232:PRO:HD3	1.82	0.60
1:X:437:ASN:HB3	7:X:891:HOH:O	2.01	0.60
1:X:327:SER:CA	7:X:915:HOH:O	2.35	0.60
1:X:226:TRP:CE2	1:X:234:SER:HB3	2.36	0.60
1:X:509:ASN:HD22	1:X:511:SER:H	1.50	0.58
1:X:194:VAL:HG11	1:X:239:VAL:HB	1.85	0.58
1:X:218:LYS:HE2	1:X:220:LYS:HE3	1.84	0.58
1:X:18:GLY:O	1:X:153:LEU:HD23	2.04	0.58
1:X:404:ILE:HD11	1:X:406:ALA:HB2	1.85	0.58
1:X:635:ILE:O	1:X:638:ILE:HG13	2.04	0.57
1:X:558:THR:HG23	1:X:587:MET:HG2	1.85	0.57
1:X:409:ASN:HD22	1:X:410:TYR:N	1.93	0.56
1:X:524:LYS:HE2	1:X:579:ASP:HB3	1.86	0.56
1:X:345:THR:CG2	1:X:443:GLU:OE2	2.53	0.56
1:X:350:MET:CE	1:X:352:LEU:HD13	2.35	0.56
1:X:403:GLN:HE22	1:X:413:SER:H	1.54	0.56
1:X:437:ASN:CB	7:X:891:HOH:O	2.53	0.56
1:X:298:THR:CG2	7:X:902:HOH:O	2.52	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:Y:73:PHE:HB3	2:Y:78:MET:SD	2.46	0.55
1:X:656:ILE:HD12	2:Y:115:VAL:CG1	2.37	0.55
1:X:345:THR:HG22	1:X:443:GLU:CD	2.27	0.54
1:X:656:ILE:HD12	2:Y:115:VAL:HG21	1.89	0.54
1:X:175:VAL:HG22	1:X:175:VAL:O	2.08	0.54
1:X:338:LEU:C	1:X:338:LEU:HD23	2.27	0.54
1:X:409:ASN:ND2	1:X:410:TYR:H	1.94	0.54
1:X:158:GLN:C	7:X:921:HOH:O	2.45	0.54
1:X:333:ALA:HB2	6:X:804:PG4:H22	1.89	0.54
2:Y:59:TRP:HB2	2:Y:114:PRO:HG2	1.90	0.54
1:X:521:MET:HE3	1:X:525:GLU:HB3	1.88	0.53
1:X:226:TRP:CZ2	1:X:234:SER:HB3	2.44	0.53
1:X:718:THR:O	1:X:721:ILE:HG13	2.09	0.53
1:X:656:ILE:HG23	2:Y:115:VAL:HG22	1.91	0.52
1:X:367:VAL:HG22	1:X:410:TYR:CD2	2.45	0.51
1:X:99:LYS:HA	1:X:99:LYS:HZ2	1.75	0.51
1:X:25:SER:O	1:X:61:GLN:HG3	2.10	0.51
1:X:699:TYR:CG	1:X:723:LYS:HD2	2.46	0.51
1:X:723:LYS:HE2	7:X:822:HOH:O	2.10	0.51
1:X:350:MET:HE2	1:X:352:LEU:HD13	1.92	0.50
1:X:69:ILE:HD11	1:X:116:ILE:HD11	1.92	0.49
1:X:74:SER:HA	1:X:108:LEU:O	2.12	0.49
1:X:17:GLN:HA	1:X:39:THR:HG22	1.94	0.49
1:X:411:TYR:C	1:X:411:TYR:CD1	2.85	0.49
2:Y:107:GLU:HB3	7:Y:419:HOH:O	2.12	0.48
1:X:153:LEU:HB3	1:X:154:PRO:HD2	1.95	0.48
1:X:240:THR:CG2	1:X:242:ARG:HD2	2.43	0.48
1:X:453:ASP:HB2	7:X:903:HOH:O	2.13	0.48
1:X:17:GLN:HE21	1:X:39:THR:HG23	1.79	0.47
1:X:203:LEU:HD23	1:X:203:LEU:C	2.35	0.47
2:Y:191:ASP:HB2	2:Y:195:GLN:HG3	1.96	0.47
1:X:61:GLN:OE1	1:X:125:PRO:HG3	2.14	0.47
1:X:585:ALA:O	1:X:586:LYS:HB2	2.15	0.47
2:Y:191:ASP:HB2	2:Y:195:GLN:CG	2.44	0.47
1:X:327:SER:CB	7:X:915:HOH:O	2.64	0.46
2:Y:206:LEU:HG	2:Y:206:LEU:O	2.15	0.46
1:X:520:ASP:HB3	7:X:914:HOH:O	2.17	0.45
2:Y:163:ALA:O	2:Y:166:SER:HB2	2.17	0.45
1:X:231:ASP:HB2	1:X:232:PRO:CD	2.47	0.45
1:X:699:TYR:CD2	1:X:723:LYS:HD2	2.52	0.45
1:X:444:LYS:HA	1:X:709:ASN:ND2	2.32	0.45
1:X:240:THR:CG2	1:X:242:ARG:HH11	2.28	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:X:593:ASP:OD1	1:X:595:ARG:CG	2.66	0.44
1:X:617:ARG:HG2	7:X:871:HOH:O	2.17	0.44
1:X:69:ILE:HG22	1:X:150:ASN:HB3	2.00	0.44
1:X:274:ASN:OD1	1:X:359:ARG:HB2	2.18	0.44
1:X:424:GLN:HG3	7:X:886:HOH:O	2.16	0.44
1:X:85:ASN:N	1:X:85:ASN:OD1	2.51	0.44
1:X:52:ASN:HB3	1:X:211:HIS:HE2	1.83	0.43
1:X:197:LYS:HB3	1:X:198:ASN:H	1.72	0.43
1:X:54:PRO:HG2	1:X:57:ASN:ND2	2.33	0.43
1:X:178:ARG:NE	1:X:224:GLU:OE2	2.49	0.43
1:X:293:THR:HB	1:X:332:VAL:CG1	2.49	0.43
1:X:360:LEU:C	1:X:360:LEU:HD23	2.40	0.43
1:X:609:GLU:O	1:X:613:LYS:HG3	2.19	0.42
1:X:298:THR:HG21	7:X:821:HOH:O	2.19	0.42
1:X:577:VAL:O	1:X:581:ILE:HG12	2.19	0.42
2:Y:98:ASP:O	2:Y:102:ILE:HD12	2.18	0.42
1:X:367:VAL:HG22	1:X:410:TYR:CE2	2.54	0.42
1:X:69:ILE:HD12	1:X:114:TYR:HB2	2.00	0.42
1:X:388:ASN:O	1:X:388:ASN:CG	2.57	0.42
1:X:460:ALA:HA	1:X:471:VAL:HA	2.01	0.42
1:X:78:THR:O	1:X:136:TRP:HA	2.19	0.42
1:X:345:THR:HG22	1:X:443:GLU:OE1	2.20	0.42
2:Y:199:VAL:CG1	2:Y:200:LYS:N	2.83	0.42
1:X:142:LYS:HE3	1:X:142:LYS:HB3	1.86	0.42
1:X:137:THR:HA	1:X:142:LYS:O	2.20	0.42
1:X:149:ASP:N	1:X:149:ASP:OD1	2.52	0.42
1:X:660:TYR:HB2	1:X:707:ILE:HG12	2.01	0.41
1:X:240:THR:HG21	1:X:242:ARG:NH1	2.31	0.41
2:Y:199:VAL:HG13	2:Y:200:LYS:N	2.35	0.41
2:Y:48:VAL:O	2:Y:145:ALA:HA	2.20	0.41
1:X:656:ILE:CD1	2:Y:115:VAL:HG11	2.47	0.41
2:Y:207:LYS:O	2:Y:207:LYS:HE2	2.21	0.41
1:X:599:ASP:HB2	7:X:876:HOH:O	2.20	0.40
2:Y:49:LEU:HD11	2:Y:66:VAL:HG21	2.02	0.40
2:Y:57:ASN:HB2	7:Y:416:HOH:O	2.21	0.40
1:X:656:ILE:HD12	2:Y:115:VAL:CG2	2.50	0.40
1:X:699:TYR:HB3	1:X:723:LYS:HB2	2.03	0.40
1:X:340:LEU:HD12	1:X:340:LEU:HA	1.89	0.40
1:X:668:LEU:HD12	1:X:668:LEU:HA	1.92	0.40

There are no symmetry-related clashes.

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	X	668/735 (91%)	630 (94%)	34 (5%)	4 (1%)	33	55
2	Y	168/189 (89%)	160 (95%)	8 (5%)	0	100	100
All	All	836/924 (90%)	790 (94%)	42 (5%)	4 (0%)	38	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	102	ASN
1	X	377	VAL
1	X	125	PRO
1	X	124	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	X	609/661 (92%)	520 (85%)	89 (15%)	5	8
2	Y	142/156 (91%)	120 (84%)	22 (16%)	4	7
All	All	751/817 (92%)	640 (85%)	111 (15%)	4	8

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	17	GLN
1	X	28	ASN
1	X	45	ILE
1	X	47	SER
1	X	55	SER

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Mol	Chain	Res	Type
1	X	66	SER
1	X	69	ILE
1	X	70	LYS
1	X	74	SER
1	X	76	GLU
1	X	93	ASP
1	X	96	VAL
1	X	99	LYS
1	X	102	ASN
1	X	105	LYS
1	X	106	ILE
1	X	107	ARG
1	X	108	LEU
1	X	110	LYS
1	X	112	ARG
1	X	115	GLN
1	X	117	LYS
1	X	122	ARG
1	X	124	ASN
1	X	127	GLU
1	X	133	LYS
1	X	136	TRP
1	X	140	GLN
1	X	149	ASP
1	X	151	LEU
1	X	197	LYS
1	X	199	LYS
1	X	207	ILE
1	X	209	ASN
1	X	212	GLU
1	X	214	LYS
1	X	216	LEU
1	X	240	THR
1	X	245	LYS
1	X	260	PRO
1	X	275	GLU
1	X	288	THR
1	X	298	THR
1	X	302	GLU
1	X	321	SER
1	X	325	SER
1	X	338	LEU

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Mol	Chain	Res	Type
1	X	340	LEU
1	X	344	ARG
1	X	345	THR
1	X	348	GLU
1	X	357	THR
1	X	388	ASN
1	X	401	LEU
1	X	404	ILE
1	X	409	ASN
1	X	413	SER
1	X	424	GLN
1	X	428	SER
1	X	437	ASN
1	X	444	LYS
1	X	447	GLN
1	X	465	GLU
1	X	466	ASN
1	X	468	ARG
1	X	470	ARG
1	X	486	GLU
1	X	509	ASN
1	X	515	GLU
1	X	518	LYS
1	X	524	LYS
1	X	527	LEU
1	X	556	GLN
1	X	557	GLN
1	X	563	LYS
1	X	576	THR
1	X	595	ARG
1	X	629	LEU
1	X	633	LYS
1	X	649	THR
1	X	654	GLU
1	X	666	SER
1	X	707	ILE
1	X	711	SER
1	X	712	GLU
1	X	713	ASN
1	X	722	LYS
1	X	723	LYS
1	X	730	LYS

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Mol	Chain	Res	Type
2	Y	51	LYS
2	Y	68	GLN
2	Y	69	LEU
2	Y	72	ARG
2	Y	74	VAL
2	Y	80	LEU
2	Y	99	ARG
2	Y	110	LYS
2	Y	111	ARG
2	Y	113	SER
2	Y	133	LYS
2	Y	137	LEU
2	Y	138	LYS
2	Y	140	SER
2	Y	160	GLU
2	Y	166	SER
2	Y	168	SER
2	Y	172	SER
2	Y	194	GLU
2	Y	195	GLN
2	Y	204	GLN
2	Y	207	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	17	GLN
1	X	28	ASN
1	X	102	ASN
1	X	104	ASN
1	X	140	GLN
1	X	158	GLN
1	X	180	ASN
1	X	209	ASN
1	X	263	HIS
1	X	268	ASN
1	X	328	ASN
1	X	363	ASN
1	X	400	GLN
1	X	403	GLN
1	X	409	ASN
1	X	466	ASN

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Mol	Chain	Res	Type
1	X	483	GLN
1	X	509	ASN
1	X	564	ASN
1	X	584	ASN
1	X	601	ASN
2	Y	64	ASN
2	Y	88	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 5 are monoatomic - leaving 1 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	$\# Z > 2$
6	PG4	X	804	-	12,12,12	0.67	0	11,11,11	0.89	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
6	PG4	X	804	-	-	0/10/10/10	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.

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 ⓘ

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	X	676/735 (91%)	0.04	27 (3%) 36 37	5, 20, 40, 49	0
2	Y	170/189 (89%)	0.25	10 (5%) 22 22	14, 23, 40, 54	0
All	All	846/924 (91%)	0.08	37 (4%) 34 34	5, 21, 40, 54	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	Y	41	ARG	5.5
1	X	124	ASN	4.9
2	Y	42	ALA	4.8
1	X	213	LYS	4.6
1	X	100	ALA	4.1
1	X	303	VAL	3.7
1	X	209	ASN	3.6
1	X	101	SER	3.4
1	X	735	GLY	3.2
2	Y	74	VAL	3.1
1	X	713	ASN	3.1
1	X	85	ASN	3.0
2	Y	208	GLY	3.0
1	X	94	GLN	2.9
1	X	514	LEU	2.8
1	X	102	ASN	2.8
1	X	670	GLN	2.7
1	X	76	GLU	2.7
1	X	320	VAL	2.6
1	X	212	GLU	2.6
1	X	140	GLN	2.5
2	Y	77	GLU	2.5
2	Y	43	PHE	2.4
1	X	425	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	X	105	LYS	2.3
2	Y	76	PRO	2.3
1	X	127	GLU	2.2
1	X	71	VAL	2.2
1	X	141	ASN	2.2
1	X	111	GLY	2.1
1	X	215	GLY	2.1
1	X	207	ILE	2.1
2	Y	138	LYS	2.1
1	X	112	ARG	2.1
2	Y	146	LEU	2.1
1	X	55	SER	2.0
2	Y	48	VAL	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	NA	X	802	1/1	0.22	4.88	28,28,28,28	0
6	PG4	X	804	13/13	0.25	4.84	27,32,37,42	0
3	CA	X	801	1/1	0.20	2.75	26,26,26,26	0
4	MN	Y	303	1/1	0.19	2.38	33,33,33,33	0
5	NA	X	803	1/1	0.14	2.12	28,28,28,28	0
3	CA	X	800	1/1	0.22	1.44	29,29,29,29	0

6.5 Other polymers ⓘ

There are no such residues in this entry.