



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:51 AM GMT

PDB ID : 2JGL
Title : CRYSTAL STRUCTURE OF MOUSE ACETYLCHOLINESTERASE INHIBITED BY AGED VX AND SARIN
Authors : Hornberg, A.; Tunemalm, A.-K.; Ekstrom, F.
Deposited on : 2007-02-13
Resolution : 2.60 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>

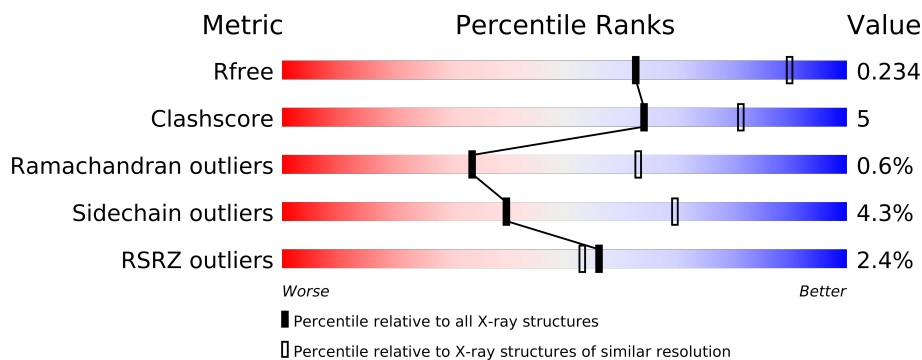
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.60 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	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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	A	548	
1	B	548	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	NAG	A	1543	-	X
3	PGE	B	1544	-	X
4	P4G	B	1545	-	X
5	P6G	B	1546	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 8712 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 ACETYLCHOLINESTERASE.

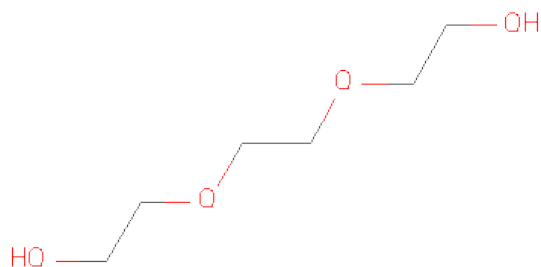
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	P	S	0	0	0
			4181	2680	725	761	1	14			
1	B	534	Total	C	N	O	P	S	0	0	1
			4163	2671	719	758	1	14			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



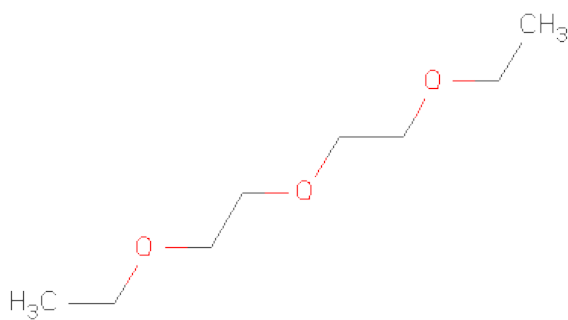
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



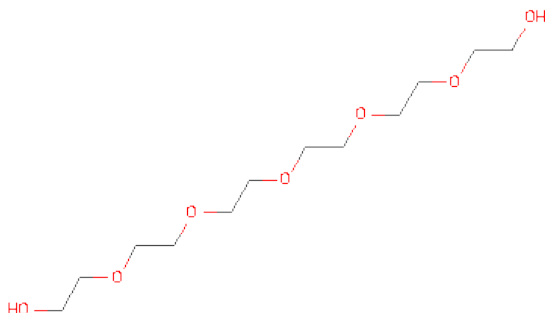
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			9	6	3		
3	B	1	Total	C	O	0	0
			9	6	3		

- Molecule 4 is 1-ETHOXY-2-(2-ETHOXYETHOXY)ETHANE (three-letter code: P4G) (formula: C₈H₁₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			11	8	3		

- Molecule 5 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $C_{12}H_{26}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			19	12	7		

- Molecule 6 is water.

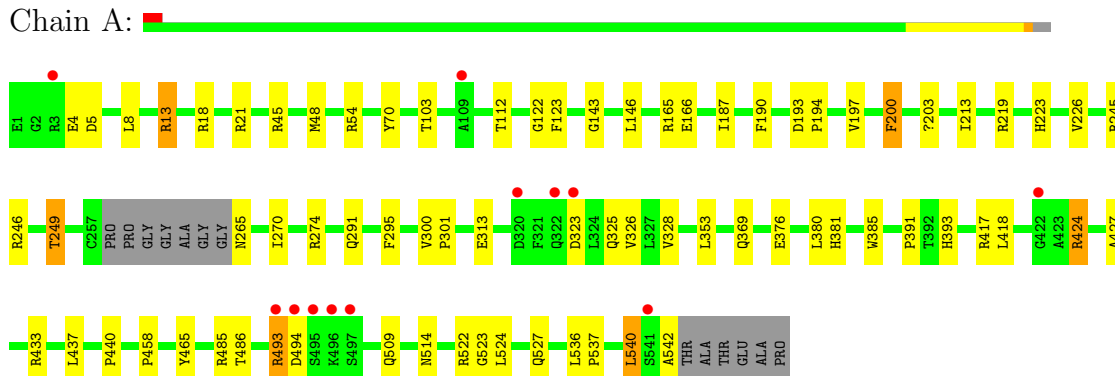
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	173	Total	O	0	0
			173	173		
6	B	133	Total	O	0	0
			133	133		

3 Residue-property plots

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.

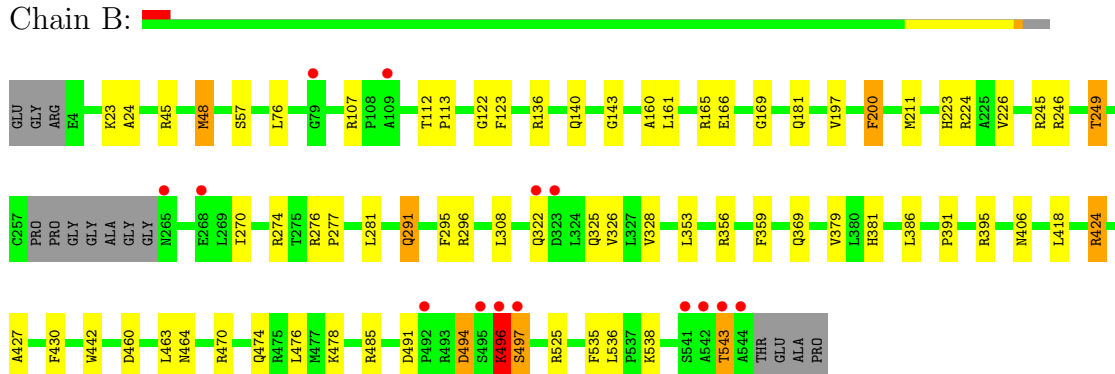
• Molecule 1: ACETYLCHOLINESTERASE

Chain A:



• Molecule 1: ACETYLCHOLINESTERASE

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.58Å 111.49Å 227.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.44 – 2.60 20.44 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.44-2.60) 100.0 (20.44-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.69 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.199 , 0.240 0.194 , 0.234	Depositor DCC
R_{free} test set	1246 reflections (2.02%)	DCC
Wilson B-factor (Å ²)	47.5	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 22.1	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 62899 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8712	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: P4G, PGE, NAG, P6G, SBG

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	0.58	0/4293	0.66	0/5864
1	B	0.58	0/4275	0.66	0/5842
All	All	0.58	0/8568	0.66	0/11706

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	A	4181	0	4066	48	1
1	B	4163	0	4049	44	1
2	A	14	0	13	0	0
3	A	9	0	11	1	0
3	B	9	0	11	0	0
4	B	11	0	18	0	0
5	B	19	0	24	3	0
6	A	173	0	0	12	0
6	B	133	0	0	8	0
All	All	8712	0	8192	90	1

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 5.

All (90) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:496:LYS:HA	6:B:2111:HOH:O	1.52	1.10
1:B:424:ARG:HH11	1:B:424:ARG:HG3	1.15	1.07
1:B:197:VAL:H	1:B:223:HIS:HD2	1.05	1.00
1:B:245:ARG:O	1:B:249:THR:HG23	1.63	0.98
1:B:197:VAL:H	1:B:223:HIS:CD2	1.93	0.87
1:B:424:ARG:HG3	1:B:424:ARG:NH1	1.84	0.81
1:A:197:VAL:H	1:A:223:HIS:HD2	1.28	0.79
1:A:48:MET:HE1	1:A:165:ARG:O	1.85	0.77
1:B:491:ASP:HB3	1:B:494:ASP:HB3	1.65	0.76
1:A:514:ASN:HB2	6:A:2161:HOH:O	1.84	0.76
1:B:424:ARG:CG	1:B:424:ARG:HH11	1.98	0.73
1:B:112:THR:HG21	1:B:143:GLY:O	1.97	0.64
1:A:197:VAL:H	1:A:223:HIS:CD2	2.16	0.63
1:A:369:GLN:HB2	6:A:2111:HOH:O	1.97	0.63
1:B:181:GLN:HG2	6:B:2054:HOH:O	1.99	0.62
1:A:380:LEU:HD23	1:A:385:TRP:HZ2	1.64	0.62
1:A:537:PRO:O	1:A:540:LEU:HB3	1.99	0.62
1:B:369:GLN:HB2	6:B:2092:HOH:O	2.00	0.62
1:A:323:ASP:HA	6:A:2120:HOH:O	1.99	0.61
1:B:381:HIS:CE1	5:B:1546:P6G:H181	2.35	0.61
1:B:460:ASP:HB3	1:B:463:LEU:HD12	1.82	0.60
1:B:296:ARG:NH2	1:B:406:ASN:OD1	2.35	0.59
1:A:265:ASN:HA	6:A:2103:HOH:O	2.02	0.59
1:B:211:MET:HG2	1:B:308:LEU:HD21	1.84	0.59
1:A:245:ARG:O	1:A:249:THR:HG22	2.03	0.58
1:A:166:GLU:HB2	1:A:270:ILE:HD13	1.88	0.56
1:A:245:ARG:O	1:A:249:THR:CG2	2.53	0.56
1:B:291:GLN:HE22	1:B:369:GLN:NE2	2.05	0.55
1:B:470:ARG:O	1:B:474:GLN:HG3	2.07	0.55
1:B:197:VAL:N	1:B:223:HIS:HD2	1.89	0.54
1:A:213:ILE:O	1:A:219:ARG:HD3	2.07	0.54
1:B:353:LEU:HB3	1:B:391:PRO:HB2	1.87	0.54
1:B:328:VAL:O	1:B:427:ALA:HA	2.08	0.54
1:B:166:GLU:HG2	1:B:270:ILE:HD13	1.93	0.50
1:B:200:PHE:CB	1:B:226:VAL:HB	2.41	0.50
1:A:393:HIS:HD2	6:A:2135:HOH:O	1.95	0.49
1:A:523:GLY:HA3	1:B:386:LEU:HD21	1.95	0.49
1:B:76:LEU:HD23	6:B:2083:HOH:O	2.13	0.48
1:B:430:PHE:HE2	1:B:476:LEU:HD11	1.79	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:540:LEU:C	1:A:542:ALA:H	2.16	0.48
1:B:326:VAL:HG21	1:B:418:LEU:HD13	1.95	0.48
3:A:1544:PGE:H6	6:A:2110:HOH:O	2.13	0.48
1:A:380:LEU:HD23	1:A:385:TRP:CZ2	2.46	0.47
1:A:313:GLU:HB2	6:A:2089:HOH:O	2.13	0.47
1:B:496:LYS:HB3	6:B:2110:HOH:O	2.14	0.47
1:A:433:ARG:CZ	1:A:437:LEU:HD23	2.44	0.47
1:A:274:ARG:HD3	6:A:2105:HOH:O	2.14	0.47
1:A:326:VAL:HG11	1:A:418:LEU:HD23	1.96	0.46
1:B:24:ALA:HB3	1:B:140:GLN:HG3	1.97	0.46
1:A:485:ARG:NH1	6:A:2156:HOH:O	2.47	0.46
1:B:291:GLN:HE21	1:B:291:GLN:HB2	1.62	0.46
1:B:496:LYS:CB	6:B:2110:HOH:O	2.64	0.46
1:A:458:PRO:HA	1:A:465:TYR:CD2	2.51	0.46
1:A:381:HIS:HA	5:B:1546:P6G:H31	1.98	0.46
1:B:160:ALA:HB2	1:B:169:GLY:CA	2.46	0.45
1:A:527:GLN:HE21	5:B:1546:P6G:H171	1.81	0.45
1:A:328:VAL:O	1:A:427:ALA:HA	2.17	0.45
1:B:166:GLU:HG2	1:B:270:ILE:CD1	2.48	0.44
1:B:224:ARG:HG2	1:B:325:GLN:HB2	1.98	0.44
1:B:359:PHE:HE1	1:B:379:VAL:HG13	1.82	0.44
1:B:200:PHE:HB2	1:B:226:VAL:HB	1.98	0.44
1:A:200:PHE:CB	1:A:226:VAL:HB	2.48	0.44
1:A:103:THR:HG21	1:A:190:PHE:HB3	1.99	0.44
1:B:113:PRO:HG2	1:B:485:ARG:HG2	2.00	0.44
1:A:5:ASP:HB3	1:A:8:LEU:HD12	2.00	0.44
1:B:395:ARG:CZ	1:B:442:TRP:HB2	2.48	0.43
1:A:48:MET:HE1	1:A:166:GLU:HA	2.01	0.43
1:A:300:VAL:HB	1:A:301:PRO:HD2	2.01	0.43
1:B:276:ARG:HA	1:B:277:PRO:HD2	1.91	0.42
1:A:219:ARG:HA	1:A:219:ARG:HD2	1.77	0.42
1:A:424:ARG:HD2	6:A:2139:HOH:O	2.19	0.42
1:A:509:GLN:HB2	6:A:2160:HOH:O	2.18	0.42
1:A:193:ASP:HA	1:A:194:PRO:HD2	1.91	0.42
1:A:353:LEU:HB3	1:A:391:PRO:HB2	2.00	0.42
1:A:323:ASP:HB2	6:A:2121:HOH:O	2.20	0.42
1:B:464:ASN:ND2	6:B:2106:HOH:O	2.47	0.42
1:A:433:ARG:NH2	1:A:440:PRO:O	2.52	0.42
1:A:112:THR:HG21	1:A:143:GLY:O	2.20	0.41
1:A:122:GLY:H	1:A:203:SBG:H1	1.68	0.41
1:A:417:ARG:HA	1:A:417:ARG:HD2	1.85	0.41
1:A:522:ARG:HE	1:A:522:ARG:HB3	1.64	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:376:GLU:CG	1:B:538:LYS:HE3	2.50	0.41
1:B:274:ARG:HD3	6:B:2073:HOH:O	2.20	0.41
1:A:122:GLY:O	1:A:123:PHE:HB2	2.21	0.41
1:A:187:ILE:HD12	1:A:187:ILE:HA	1.95	0.41
1:A:4:GLU:OE2	1:A:18:ARG:HD3	2.21	0.41
1:A:485:ARG:HB3	1:A:486:THR:HG23	2.03	0.41
1:B:48:MET:HE1	1:B:165:ARG:O	2.21	0.41
1:A:380:LEU:HD13	1:B:535:PHE:HB2	2.02	0.41
1:B:122:GLY:O	1:B:123:PHE: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	Distance(Å)	Clash(Å)
1:A:13:ARG:NH1	1:B:57:SER:O[2.555]	2.03	0.17

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	530/548 (97%)	501 (94%)	26 (5%)	3 (1%)	33	63
1	B	529/548 (96%)	507 (96%)	19 (4%)	3 (1%)	33	63
All	All	1059/1096 (97%)	1008 (95%)	45 (4%)	6 (1%)	33	63

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	494	ASP
1	B	496	LYS
1	B	543	THR
1	B	497	SER
1	A	493	ARG
1	A	540	LEU

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	438/445 (98%)	422 (96%)	16 (4%)	45	75
1	B	437/445 (98%)	415 (95%)	22 (5%)	34	61
All	All	875/890 (98%)	837 (96%)	38 (4%)	40	69

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	21	ARG
1	A	45	ARG
1	A	54	ARG
1	A	70	TYR
1	A	146	LEU
1	A	200	PHE
1	A	246	ARG
1	A	249	THR
1	A	291	GLN
1	A	295	PHE
1	A	325	GLN
1	A	424	ARG
1	A	493	ARG
1	A	524	LEU
1	A	536	LEU
1	B	23	LYS
1	B	45	ARG
1	B	48	MET
1	B	107	ARG
1	B	136	ARG
1	B	161	LEU
1	B	200	PHE
1	B	246	ARG
1	B	249	THR
1	B	281	LEU
1	B	291	GLN
1	B	295	PHE
1	B	322	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	356	ARG
1	B	424	ARG
1	B	478	LYS
1	B	494	ASP
1	B	496	LYS
1	B	497	SER
1	B	525	ARG
1	B	536	LEU
1	B	543	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	HIS
1	A	287	HIS
1	A	291	GLN
1	B	223	HIS
1	B	291	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

2 non-standard protein/DNA/RNA residues are 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	$\# Z > 2$
1	SBG	A	203	1	9,9,10	6.74	3 (33%)	10,12,14	3.09	4 (40%)
1	SBG	B	203	1	9,9,10	6.44	2 (22%)	10,12,14	2.96	5 (50%)

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
1	SBG	A	203	1	-	0/6/8/10	0/0/0/0
1	SBG	B	203	1	-	0/6/8/10	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	203	SBG	O-C	19.38	1.24	1.11
1	B	203	SBG	O-C	18.72	1.24	1.11
1	A	203	SBG	CA-C	4.94	1.57	1.48
1	B	203	SBG	CA-C	4.16	1.56	1.48
1	A	203	SBG	P1-OG	2.48	1.61	1.58

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	203	SBG	P1-OG-CB	-8.54	110.64	122.10
1	B	203	SBG	P1-OG-CB	-6.69	113.12	122.10
1	B	203	SBG	O1-P1-O2	3.41	119.69	110.68
1	B	203	SBG	OG-P1-C1	-3.40	87.74	105.23
1	B	203	SBG	OG-P1-O2	3.31	122.26	111.99
1	A	203	SBG	O1-P1-O2	2.76	117.95	110.68
1	A	203	SBG	OG-P1-C1	-2.59	91.90	105.23
1	A	203	SBG	OG-CB-CA	2.50	112.23	108.69
1	B	203	SBG	OG-CB-CA	2.46	112.17	108.69

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

5 ligands are 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	NAG	A	1543	1	12,14,15	0.64	0	15,19,21	1.54	2 (13%)
3	PGE	A	1544	-	8,8,9	1.59	1 (12%)	6,7,8	0.29	0
3	PGE	B	1544	-	8,8,9	11.72	2 (25%)	6,7,8	0.48	0
4	P4G	B	1545	-	10,10,10	0.65	0	9,9,9	0.32	0
5	P6G	B	1546	-	18,18,18	1.69	5 (27%)	17,17,17	2.09	9 (52%)

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
2	NAG	A	1543	1	-	0/6/23/26	0/1/1/1
3	PGE	A	1544	-	-	0/6/6/7	0/0/0/0
3	PGE	B	1544	-	-	0/6/6/7	0/0/0/0
4	P4G	B	1545	-	-	0/8/8/8	0/0/0/0
5	P6G	B	1546	-	-	0/16/16/16	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1544	PGE	C1-C2	32.99	1.68	1.55
3	A	1544	PGE	C1-C2	-4.29	1.53	1.55
5	B	1546	P6G	C9-C8	-2.92	1.33	1.48
5	B	1546	P6G	C3-C2	-2.77	1.34	1.49
5	B	1546	P6G	C15-C14	-2.76	1.34	1.48
3	B	1544	PGE	O2-C2	2.75	1.52	1.42
5	B	1546	P6G	C17-C18	-2.62	1.35	1.49
5	B	1546	P6G	C6-C5	-2.54	1.35	1.48

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1543	NAG	O5-C5-C6	3.77	110.94	106.98
5	B	1546	P6G	O7-C6-C5	3.14	124.72	110.47
5	B	1546	P6G	O4-C5-C6	2.96	123.91	110.47
5	B	1546	P6G	O13-C14-C15	2.78	123.08	110.47
5	B	1546	P6G	O7-C8-C9	2.73	122.87	110.47
5	B	1546	P6G	O16-C15-C14	2.61	122.30	110.47
2	A	1543	NAG	O3-C3-C2	2.33	113.98	109.09
5	B	1546	P6G	C17-O16-C15	2.27	123.35	113.38
5	B	1546	P6G	O4-C3-C2	2.15	120.93	110.61
5	B	1546	P6G	O10-C9-C8	2.11	120.05	110.47
5	B	1546	P6G	C5-O4-C3	2.02	122.27	113.38

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	A	535/548 (97%)	-0.43	12 (2%) 59 56	21, 37, 56, 80	0
1	B	534/548 (97%)	-0.30	14 (2%) 53 50	27, 42, 62, 85	0
All	All	1069/1096 (97%)	-0.36	26 (2%) 56 53	21, 39, 60, 85	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	544	ALA	7.7
1	B	497	SER	5.1
1	B	541	SER	4.5
1	B	543	THR	4.5
1	A	493	ARG	4.0
1	A	494	ASP	3.9
1	B	542	ALA	3.8
1	B	268	GLU	3.4
1	A	495	SER	3.4
1	A	496	LYS	3.4
1	A	322	GLN	3.1
1	A	323	ASP	3.0
1	B	323	ASP	3.0
1	A	541	SER	2.8
1	A	497	SER	2.7
1	B	495	SER	2.7
1	B	109	ALA	2.6
1	A	422	GLY	2.4
1	B	322	GLN	2.3
1	A	109	ALA	2.3
1	B	496	LYS	2.2
1	A	3	ARG	2.2
1	B	492	PRO	2.1
1	A	320	ASP	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	265	ASN	2.1
1	B	79	GLY	2.1

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

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
1	SBG	B	203	10/11	0.12	0.46	31,32,35,35	0
1	SBG	A	203	10/11	0.09	-0.26	25,27,29,30	0

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
2	NAG	A	1543	14/15	0.39	11.66	66,71,77,78	0
4	P4G	B	1545	11/11	0.23	7.87	87,89,90,91	0
3	PGE	B	1544	9/10	0.32	3.06	81,84,85,86	0
5	P6G	B	1546	19/19	0.25	2.22	53,62,67,68	0
3	PGE	A	1544	9/10	0.25	1.52	71,74,77,78	0

6.5 Other polymers ⓘ

There are no such residues in this entry.