



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 05:25 am GMT

PDB ID : 2ACH
Title : CRYSTAL STRUCTURE OF CLEAVED HUMAN ALPHA1-ANTICHYMO
TRYPSIN AT 2.7 ANGSTROMS RESOLUTION AND ITS COMPARISON
WITH OTHER SERPINS
Authors : Baumann, U.; Huber, R.; Bode, W.; Grosse, D.; Lesjak, M.; Laurell, C.B.
Deposited on : 1993-04-26
Resolution : 2.70 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

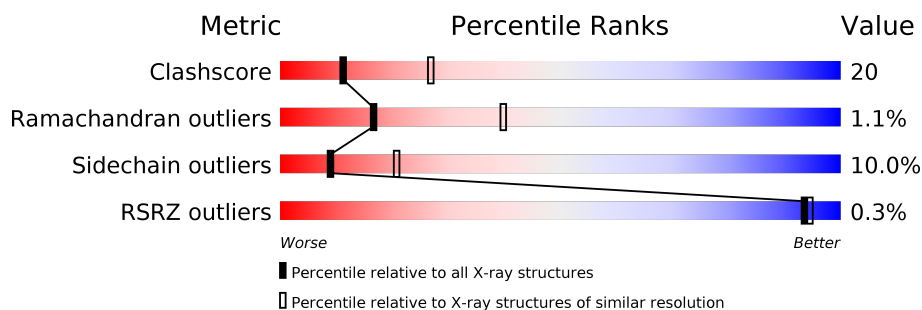
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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 $\leq 5\%$. 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	360	
2	B	40	

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
4	NAG	A	540	-	-	-	X
5	PO4	A	550	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3126 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 ALPHA 1-ANTICHYMOTRYPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	40	0	0
			2678	1718	437	512	11			

- Molecule 2 is a protein called ALPHA 1-ANTICHYMOTRYPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	35	Total	C	N	O	S	13	0	0
			290	188	51	49	2			

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	3	0
			28	16	2	10		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

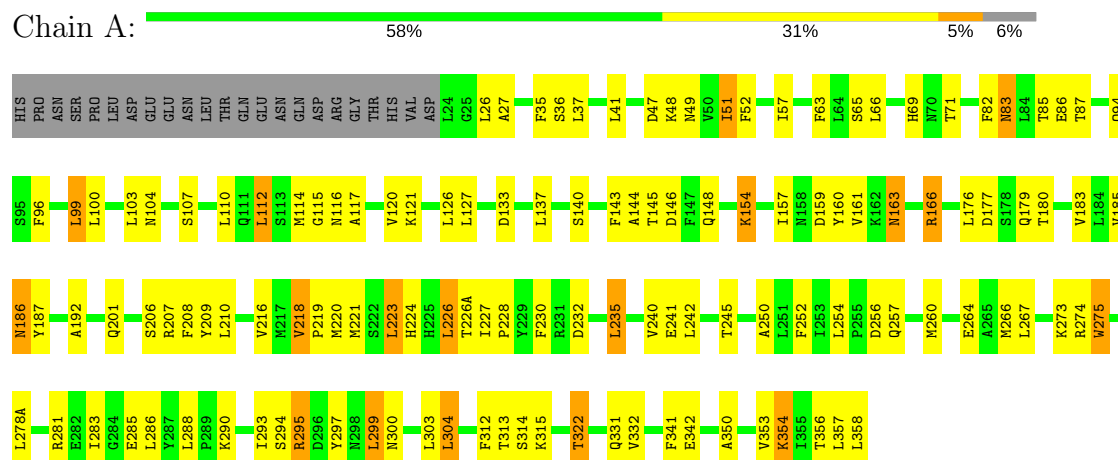
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	80	Total	O	0	0
			80	80		
6	B	6	Total	O	0	0
			6	6		

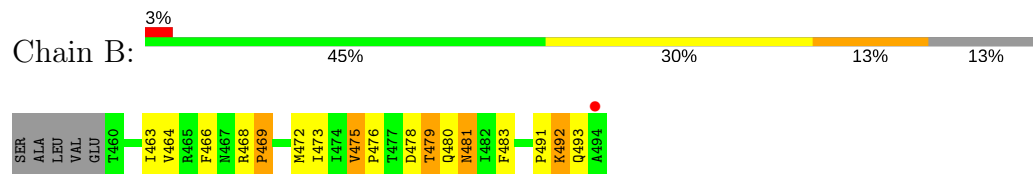
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: ALPHA 1-ANTICHYMOTRYPSIN



• Molecule 2: ALPHA 1-ANTICHYMOTRYPSIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	75.30Å 75.30Å 208.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.70 26.17 – 2.66	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.70) 77.3 (26.17-2.66)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.181 , (Not available) 0.187 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 126.4	EDS
L-test for twinning ¹	$\langle L \rangle = 0.32$, $\langle L^2 \rangle = 0.15$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3126	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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: PO4, BMA, NAG

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	0.53	0/2728	0.76	0/3686
2	B	0.55	0/296	0.89	0/399
All	All	0.53	0/3024	0.77	0/4085

There are no bond length outliers.

There are no bond angle outliers.

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	2678	0	2694	107	0
2	B	290	0	304	26	0
3	A	39	0	34	0	0
4	A	28	0	25	1	0
5	A	5	0	0	2	0
6	A	80	0	0	3	0
6	B	6	0	0	1	0
All	All	3126	0	3057	119	0

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 20.

All (119) 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:303:LEU:HD12	1:A:312:PHE:HZ	1.41	0.86
1:A:110:LEU:HD13	1:A:192:ALA:HB2	1.56	0.85
1:A:230:PHE:CZ	1:A:232:ASP:HB3	2.18	0.78
2:B:478:ASP:O	2:B:479:THR:HB	1.81	0.78
1:A:254:LEU:HB2	2:B:466:PHE:CE2	2.20	0.75
1:A:300:ASN:O	1:A:304:LEU:HD12	1.91	0.71
1:A:37:LEU:HD23	1:A:52:PHE:CE2	2.26	0.70
1:A:285:GLU:HB2	2:B:463:ILE:HG12	1.73	0.69
1:A:256:ASP:HA	2:B:468:ARG:HH21	1.59	0.68
1:A:69:HIS:CD2	5:A:550:PO4:O4	2.48	0.66
1:A:117:ALA:HB1	1:A:143:PHE:HE2	1.60	0.66
1:A:110:LEU:CD1	1:A:192:ALA:HB2	2.26	0.66
1:A:177:ASP:HB3	1:A:180:THR:OG1	1.95	0.66
1:A:83:ASN:HD21	1:A:85:THR:HB	1.62	0.64
1:A:35:PHE:CE1	1:A:267:LEU:HD22	2.34	0.63
1:A:303:LEU:HD12	1:A:312:PHE:CZ	2.30	0.63
1:A:154:LYS:HG3	1:A:176:LEU:HB2	1.81	0.62
1:A:242:LEU:HD12	1:A:252:PHE:CE1	2.35	0.62
1:A:206:SER:HB2	1:A:220:MET:HB2	1.81	0.62
1:A:226(A):THR:HA	1:A:281:ARG:O	2.01	0.61
1:A:37:LEU:HD23	1:A:52:PHE:HE2	1.64	0.61
2:B:475:VAL:HG23	2:B:476:PRO:HD2	1.83	0.60
1:A:313:THR:HG22	1:A:315:LYS:H	1.66	0.59
1:A:300:ASN:HD22	1:A:300:ASN:N	2.01	0.58
1:A:41:LEU:HD22	1:A:297:TYR:CE1	2.38	0.58
1:A:121:LYS:HA	1:A:145:THR:O	2.03	0.58
1:A:235:LEU:HD12	1:A:274:ARG:HH22	1.69	0.58
1:A:83:ASN:ND2	1:A:86:GLU:HG2	2.18	0.58
1:A:127:LEU:HG	1:A:322:THR:O	2.04	0.57
1:A:313:THR:HG22	1:A:314:SER:N	2.19	0.56
1:A:223:LEU:HD23	1:A:224:HIS:H	1.69	0.56
2:B:476:PRO:HD2	2:B:481:ASN:O	2.05	0.56
1:A:226:LEU:HD12	1:A:283:ILE:HB	1.88	0.55
1:A:117:ALA:HB1	1:A:143:PHE:CE2	2.42	0.55
1:A:226:LEU:CD1	1:A:283:ILE:HB	2.37	0.55
1:A:300:ASN:HD21	1:A:332:VAL:H	1.55	0.54
1:A:356:THR:HG22	1:A:357:LEU:N	2.23	0.54
2:B:480:GLN:HB3	6:B:617:HOH:O	2.06	0.54
1:A:185:VAL:HG12	1:A:186:ASN:N	2.22	0.54
1:A:207:ARG:HA	1:A:216:VAL:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:492:LYS:HD3	2:B:492:LYS:O	2.08	0.53
1:A:295:ARG:HH21	2:B:493:GLN:NE2	2.05	0.53
1:A:314:SER:HA	1:A:358:LEU:CD2	2.38	0.53
1:A:186:ASN:C	1:A:186:ASN:ND2	2.63	0.52
1:A:235:LEU:HD13	1:A:275:TRP:HH2	1.74	0.52
1:A:48:LYS:HB3	1:A:295:ARG:HH12	1.75	0.52
2:B:476:PRO:HB2	2:B:478:ASP:OD1	2.10	0.51
1:A:110:LEU:HD13	1:A:192:ALA:CB	2.36	0.51
1:A:299:LEU:O	1:A:303:LEU:HG	2.10	0.51
1:A:100:LEU:O	1:A:104:ASN:HB2	2.09	0.51
1:A:27:ALA:HB1	2:B:481:ASN:ND2	2.26	0.51
1:A:83:ASN:ND2	1:A:85:THR:HB	2.25	0.51
1:A:159:ASP:O	1:A:163:ASN:HB2	2.10	0.51
1:A:275:TRP:O	1:A:278(A):LEU:HG	2.10	0.51
1:A:69:HIS:HD2	5:A:550:PO4:O4	1.94	0.51
1:A:223:LEU:HD13	1:A:226:LEU:HD23	1.93	0.50
2:B:479:THR:CG2	2:B:480:GLN:N	2.75	0.50
1:A:120:VAL:HG11	1:A:126:LEU:HD21	1.93	0.50
1:A:208:PHE:O	1:A:216:VAL:HG12	2.11	0.50
1:A:144:ALA:HB2	6:A:668:HOH:O	2.12	0.49
1:A:230:PHE:CE2	1:A:232:ASP:HB3	2.47	0.49
2:B:478:ASP:O	2:B:479:THR:CB	2.55	0.48
1:A:286:LEU:HA	2:B:464:VAL:O	2.14	0.48
1:A:186:ASN:O	1:A:350:ALA:HA	2.14	0.48
1:A:37:LEU:HD21	1:A:57:ILE:HD13	1.96	0.48
1:A:290:LYS:HG3	1:A:341:PHE:HA	1.96	0.47
1:A:115:GLY:HA2	4:A:540:NAG:O7	2.14	0.47
1:A:295:ARG:HH21	2:B:493:GLN:HE21	1.62	0.47
1:A:96:PHE:O	1:A:100:LEU:HG	2.14	0.47
1:A:260:MET:HG3	1:A:264:GLU:OE2	2.14	0.47
1:A:99:LEU:HD23	1:A:100:LEU:N	2.29	0.47
1:A:235:LEU:HD13	1:A:275:TRP:CH2	2.49	0.47
1:A:232:ASP:OD2	1:A:235:LEU:HB2	2.16	0.46
1:A:112:LEU:HD21	2:B:483:PHE:CE2	2.50	0.46
1:A:157:ILE:O	1:A:161:VAL:HG23	2.15	0.46
1:A:295:ARG:NH2	2:B:493:GLN:HE21	2.14	0.46
1:A:117:ALA:HB2	1:A:160:TYR:CZ	2.50	0.46
1:A:227:ILE:HD12	1:A:240:VAL:HG13	1.98	0.45
1:A:313:THR:CG2	1:A:314:SER:N	2.79	0.45
1:A:120:VAL:O	1:A:145:THR:N	2.47	0.45
1:A:186:ASN:HD22	1:A:186:ASN:C	2.19	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:473:ILE:HG22	2:B:475:VAL:HG12	1.97	0.45
1:A:331:GLN:NE2	1:A:354:LYS:NZ	2.64	0.45
1:A:218:VAL:HG22	1:A:219:PRO:HD2	1.99	0.45
1:A:145:THR:HG22	1:A:146:ASP:N	2.32	0.44
1:A:103:LEU:HB2	1:A:114:MET:CE	2.47	0.44
1:A:230:PHE:CB	1:A:278(A):LEU:HD23	2.48	0.44
1:A:63:PHE:O	1:A:66:LEU:HB3	2.18	0.44
1:A:218:VAL:HG21	2:B:491:PRO:HB2	1.98	0.44
1:A:230:PHE:HB2	1:A:278(A):LEU:HD23	2.00	0.44
1:A:235:LEU:HD21	1:A:266:MET:HB2	1.99	0.44
1:A:37:LEU:HD21	1:A:57:ILE:CD1	2.48	0.44
1:A:133:ASP:O	1:A:137:LEU:HB2	2.18	0.43
1:A:146:ASP:C	1:A:148:GLN:H	2.21	0.43
1:A:223:LEU:HD23	1:A:224:HIS:N	2.32	0.43
1:A:224:HIS:CD2	1:A:285:GLU:HG3	2.52	0.43
1:A:228:PRO:HB2	1:A:241:GLU:HB3	2.00	0.43
1:A:257:GLN:NE2	6:A:637:HOH:O	2.51	0.43
1:A:221:MET:CG	1:A:290:LYS:HB2	2.48	0.43
1:A:250:ALA:HB1	2:B:472:MET:HE3	2.00	0.43
2:B:475:VAL:HG23	2:B:476:PRO:CD	2.47	0.43
1:A:313:THR:HG21	1:A:315:LYS:HE2	2.00	0.42
1:A:209:TYR:O	2:B:469:PRO:HD3	2.20	0.42
1:A:183:VAL:HA	1:A:353:VAL:O	2.19	0.42
1:A:166:ARG:NH1	6:A:615:HOH:O	2.52	0.42
1:A:283:ILE:HD12	2:B:464:VAL:HG23	2.02	0.42
1:A:115:GLY:HA3	1:A:187:TYR:CE2	2.55	0.41
1:A:300:ASN:HD21	1:A:332:VAL:N	2.18	0.41
1:A:47:ASP:OD1	1:A:47:ASP:N	2.53	0.41
1:A:300:ASN:ND2	1:A:300:ASN:N	2.67	0.41
1:A:254:LEU:HB2	2:B:466:PHE:HE2	1.81	0.41
1:A:26:LEU:HD22	1:A:82:PHE:HD1	1.86	0.41
1:A:331:GLN:HB3	1:A:354:LYS:HD3	2.02	0.40
2:B:473:ILE:HG22	2:B:473:ILE:O	2.22	0.40
1:A:221:MET:HB2	1:A:288:LEU:O	2.21	0.40
1:A:51:ILE:HG21	1:A:293:ILE:HG13	2.03	0.40
1:A:116:ASN:O	1:A:140:SER:HA	2.22	0.40
1:A:154:LYS:HE2	1:A:154:LYS:HB3	1.85	0.40
2:B:468:ARG:HH11	2:B:468:ARG:HG2	1.86	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	A	335/360 (93%)	301 (90%)	31 (9%)	3 (1%)	20	46
2	B	33/40 (82%)	28 (85%)	4 (12%)	1 (3%)	5	12
All	All	368/400 (92%)	329 (89%)	35 (10%)	4 (1%)	17	40

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	479	THR
1	A	107	SER
1	A	154	LYS
1	A	51	ILE

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	296/318 (93%)	267 (90%)	29 (10%)	9	21
2	B	34/38 (90%)	30 (88%)	4 (12%)	6	14
All	All	330/356 (93%)	297 (90%)	33 (10%)	9	21

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	SER
1	A	49	ASN

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Mol	Chain	Res	Type
1	A	65	SER
1	A	71	THR
1	A	83	ASN
1	A	87	THR
1	A	94	GLN
1	A	99	LEU
1	A	112	LEU
1	A	163	ASN
1	A	166	ARG
1	A	179	GLN
1	A	186	ASN
1	A	201	GLN
1	A	210	LEU
1	A	218	VAL
1	A	223	LEU
1	A	226	LEU
1	A	235	LEU
1	A	245	THR
1	A	273	LYS
1	A	275	TRP
1	A	294	SER
1	A	295	ARG
1	A	299	LEU
1	A	304	LEU
1	A	322	THR
1	A	342	GLU
1	A	354	LYS
2	B	469	PRO
2	B	475	VAL
2	B	481	ASN
2	B	492	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	HIS
1	A	83	ASN
1	A	163	ASN
1	A	186	ASN
1	A	201	GLN
1	A	204	HIS
1	A	300	ASN

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Mol	Chain	Res	Type
1	A	331	GLN
2	B	481	ASN
2	B	493	GLN

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 ⓘ

5 carbohydrates 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
3	NAG	A	520	1,3	14,14,15	0.55	0	15,19,21	0.98	1 (6%)
3	NAG	A	522	3	14,14,15	0.96	1 (7%)	15,19,21	1.70	4 (26%)
3	BMA	A	523	3	11,11,12	0.53	0	13,15,17	0.97	1 (7%)
4	NAG	A	540	1,4	14,14,15	0.76	0	15,19,21	1.56	2 (13%)
4	NAG	A	542	4	14,14,15	0.67	0	15,19,21	1.26	1 (6%)

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	NAG	A	520	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	522	3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BMA	A	523	3	-	0/2/19/22	0/1/1/1
4	NAG	A	540	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	542	4	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	522	NAG	C4-C3	2.54	1.58	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	540	NAG	C4-C3-C2	-4.37	104.61	111.02
4	A	542	NAG	C1-C2-N2	-3.73	104.11	110.49
3	A	522	NAG	C4-C3-C2	-3.19	106.34	111.02
3	A	520	NAG	C1-C2-N2	-2.47	106.27	110.49
3	A	522	NAG	O3-C3-C2	-2.31	104.43	109.39
3	A	523	BMA	C1-C2-C3	-2.11	106.98	109.65
3	A	522	NAG	O3-C3-C4	2.04	114.80	110.36
4	A	540	NAG	O4-C4-C3	2.44	115.67	110.36
3	A	522	NAG	O4-C4-C3	4.29	119.70	110.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	540	NAG	1	0

5.6 Ligand geometry

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
5	PO4	A	550	-	4,4,4	0.86	0	6,6,6	0.39	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
5	PO4	A	550	-	-	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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	550	PO4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	337/360 (93%)	-0.87	0	100 100	3, 20, 45, 59	10 (2%)
2	B	35/40 (87%)	-0.41	1 (2%)	52 52	5, 16, 55, 61	4 (11%)
All	All	372/400 (93%)	-0.83	1 (0%)	93 94	3, 20, 48, 61	14 (3%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	494	ALA	3.1

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](#)

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	NAG	A	540	14/15	0.93	0.15	2.50	38,46,59,59	0
3	NAG	A	520	14/15	0.95	0.12	0.39	27,30,37,38	0
4	NAG	A	542	14/15	0.91	0.21	-	61,65,72,72	3
3	NAG	A	522	14/15	0.94	0.23	-	27,45,54,57	0
3	BMA	A	523	11/12	0.82	0.21	-	65,69,72,75	0

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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	PO4	A	550	5/5	0.99	0.13	-	22,22,22,22	5

6.5 Other polymers [i](#)

There are no such residues in this entry.