Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 01:40 AM GMT

PDB ID	:	2DW2
Title	:	Crystal structure of VAP2 from Crotalus atrox venom (Form 2-5 crystal)
Authors	:	Takeda, S.; Igarashi, T.; Araki, S.
Deposited on	:	2006-08-02
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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

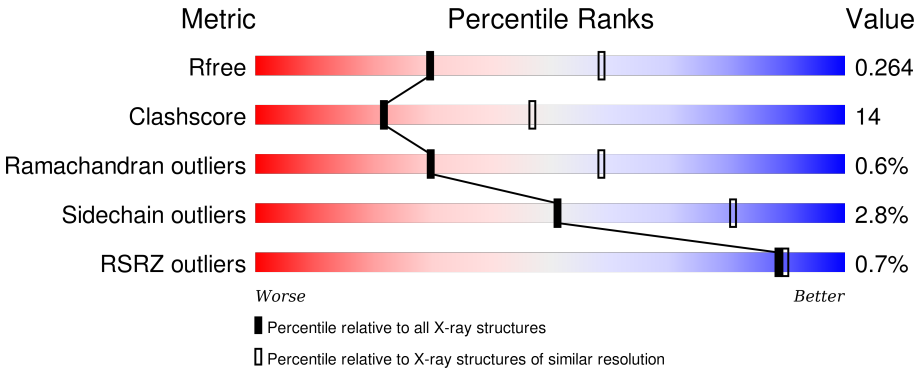
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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))
R_{free}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (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 <=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	419	% 74% 24%
1	B	419	% 70% 27%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6823 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 Catrocollastatin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	415	Total	C	N	O	S	0	0	0
			3219	1990	545	637	47			
1	B	415	Total	C	N	O	S	0	0	0
			3219	1990	545	637	47			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	203	VAL	PHE	SEE REMARK 999	UNP Q90282
B	203	VAL	PHE	SEE REMARK 999	UNP Q90282

- Molecule 2 is a polymer of unknown type called SUGAR (9-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	9	Total	C	N	O	0	0
			113	64	5	44		
2	B	9	Total	C	N	O	0	0
			113	64	5	44		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	3	Total	Ca	0	0
			3	3		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	3	Total Ca 3 3	0	0

- Molecule 5 is water.

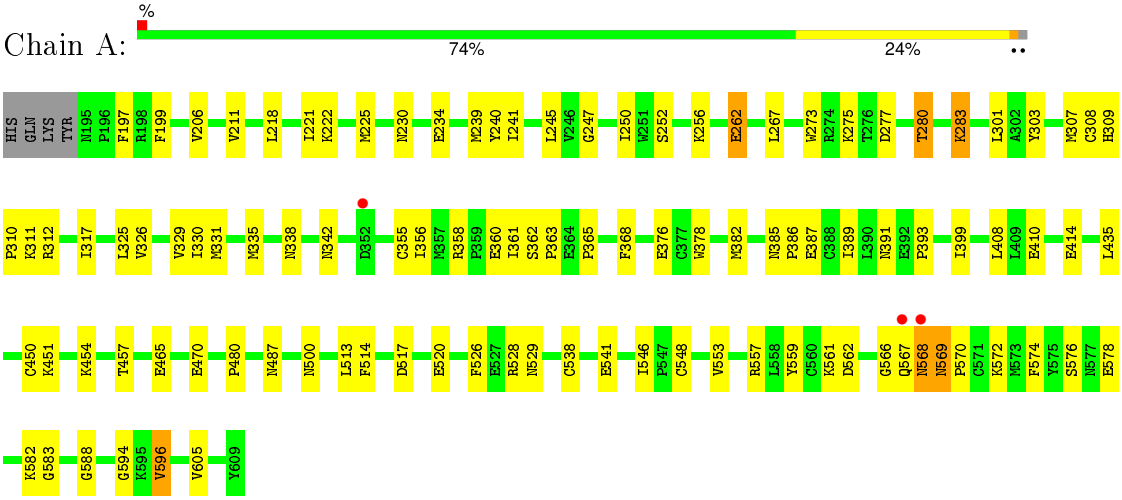
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	85	Total O 85 85	0	0
5	B	66	Total O 66 66	0	0

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.

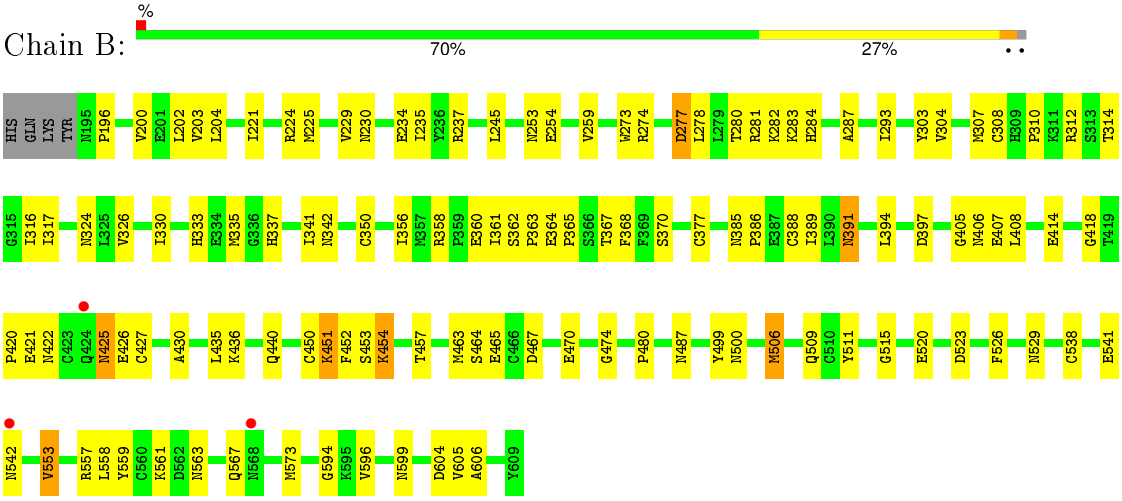
- Molecule 1: Catrocollastatin

Chain A:



- Molecule 1: Catrocollastatin

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	220.74Å 79.47Å 58.69Å 90.00° 91.73° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 46.38 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.8 (50.00-2.70) 95.6 (46.38-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.31 (at 2.69Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.199 , 0.260 0.214 , 0.264	Depositor DCC
R_{free} test set	1334 reflections (4.96%)	DCC
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 46.1	EDS
Estimated twinning fraction	0.094 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 26911 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6823	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN, BMA, NAG, CA, FUC, MAN

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	0.29	0/3291	0.55	0/4453
1	B	0.28	0/3291	0.55	0/4453
All	All	0.28	0/6582	0.55	0/8906

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	3219	0	2985	82	0
1	B	3219	0	2985	98	0
2	A	113	0	97	5	0
2	B	113	0	97	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
5	A	85	0	0	4	0
5	B	66	0	0	5	0
All	All	6823	0	6164	180	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 14.

All (180) 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:529:ASN:HD21	1:A:557:ARG:HB3	1.36	0.91
1:B:326:VAL:HG13	1:B:361:ILE:HD11	1.63	0.81
1:B:529:ASN:HD21	1:B:557:ARG:HB3	1.47	0.80
1:B:454:LYS:HE3	1:B:457:THR:HG21	1.69	0.72
1:B:454:LYS:CD	1:B:454:LYS:H	2.02	0.72
1:B:224:ARG:HE	1:B:324:ASN:ND2	1.87	0.71
1:B:454:LYS:H	1:B:454:LYS:HD3	1.53	0.71
1:A:470:GLU:HB3	1:A:480:PRO:HG2	1.75	0.69
1:A:378:TRP:O	1:A:382:MET:HG2	1.93	0.69
1:B:224:ARG:HE	1:B:324:ASN:HD21	1.42	0.68
1:A:574:PHE:CZ	1:A:582:LYS:HE3	2.29	0.68
1:B:526:PHE:O	1:B:553:VAL:HG22	1.94	0.67
1:B:386:PRO:O	1:B:389:ILE:HG12	1.94	0.67
1:A:275:LYS:HE2	1:A:311:LYS:HB2	1.77	0.67
1:A:526:PHE:HB3	1:A:553:VAL:HG13	1.78	0.66
1:A:326:VAL:HG13	1:A:361:ILE:HD11	1.78	0.65
1:B:425:ASN:HD22	1:B:426:GLU:N	1.95	0.64
1:A:408:LEU:O	1:A:410:GLU:HG3	1.96	0.64
1:B:363:PRO:C	1:B:365:PRO:HD3	2.19	0.63
1:A:277:ASP:O	1:A:280:THR:HG22	1.98	0.63
1:A:526:PHE:HB3	1:A:553:VAL:CG1	2.29	0.63
1:B:277:ASP:O	1:B:280:THR:HB	1.98	0.63
1:B:326:VAL:HG13	1:B:361:ILE:CD1	2.29	0.62
1:A:562:ASP:CG	1:A:567:GLN:HG3	2.20	0.61
1:A:355:CYS:H	1:A:358:ARG:NH1	1.98	0.61
2:A:806:NAG:H3	1:B:465:GLU:OE2	2.01	0.61
1:A:317:ILE:HD13	1:A:331:MET:HA	1.82	0.61
1:B:204:LEU:HD13	1:B:225:MET:CE	2.30	0.61
1:A:317:ILE:CD1	1:A:331:MET:HA	2.30	0.60
1:B:594:GLY:C	1:B:605:VAL:HG23	2.21	0.60
1:B:283:LYS:HD2	1:B:283:LYS:H	1.66	0.60
1:A:487:ASN:ND2	1:A:500:ASN:H	2.00	0.60
1:B:558:LEU:HD12	1:B:559:TYR:H	1.67	0.59
1:B:204:LEU:HD13	1:B:225:MET:HE3	1.84	0.59
1:B:203:VAL:HG11	1:B:278:LEU:HD21	1.83	0.59
1:A:206:VAL:HB	1:A:250:ILE:HD13	1.83	0.59
1:B:350:CYS:HB3	1:B:370:SER:HA	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:517:ASP:OD2	1:B:342:ASN:HB3	2.03	0.59
1:B:391:ASN:HD22	1:B:391:ASN:N	2.00	0.59
1:A:568:ASN:O	1:A:570:PRO:HD3	2.04	0.58
1:A:487:ASN:HD21	1:A:500:ASN:H	1.51	0.58
1:B:229:VAL:HG21	1:B:245:LEU:HD13	1.86	0.58
1:B:596:VAL:HG13	1:B:605:VAL:HG22	1.86	0.57
1:B:283:LYS:N	1:B:283:LYS:HD2	2.19	0.57
1:B:454:LYS:HD3	1:B:454:LYS:N	2.20	0.57
1:B:361:ILE:HB	5:B:27:HOH:O	2.04	0.57
1:B:440:GLN:HE21	1:B:452:PHE:H	1.53	0.57
1:A:283:LYS:N	1:A:283:LYS:HD2	2.20	0.56
1:B:487:ASN:HD21	1:B:500:ASN:H	1.53	0.56
1:A:454:LYS:O	1:A:457:THR:HG23	2.04	0.56
1:A:567:GLN:O	1:A:568:ASN:HB3	2.05	0.56
1:A:326:VAL:HG13	1:A:361:ILE:CD1	2.35	0.56
1:A:312:ARG:HD3	1:B:567:GLN:OE1	2.05	0.56
1:B:425:ASN:ND2	1:B:427:CYS:H	2.03	0.56
1:A:245:LEU:HG	1:A:399:ILE:HD11	1.88	0.55
1:A:307:MET:O	1:A:308:CYS:HB2	2.06	0.55
1:B:229:VAL:HG11	1:B:245:LEU:HB2	1.89	0.54
1:B:561:LYS:HD3	1:B:563:ASN:HD21	1.72	0.54
1:B:454:LYS:H	1:B:454:LYS:CE	2.21	0.54
1:A:221:ILE:O	1:A:225:MET:HG2	2.08	0.54
2:A:804:NAG:H61	2:A:806:NAG:C8	2.37	0.54
1:B:304:VAL:HA	1:B:337:HIS:O	2.08	0.54
1:B:440:GLN:NE2	1:B:452:PHE:H	2.05	0.53
1:B:511:TYR:HA	1:B:515:GLY:O	2.08	0.53
1:A:360:GLU:HG2	5:A:67:HOH:O	2.08	0.53
1:B:221:ILE:O	1:B:225:MET:HG2	2.08	0.53
1:B:307:MET:O	1:B:308:CYS:HB2	2.09	0.53
1:B:259:VAL:HG12	1:B:293:ILE:HD12	1.91	0.53
1:B:364:GLU:N	1:B:365:PRO:HD3	2.23	0.53
1:B:450:CYS:C	1:B:451:LYS:HD2	2.29	0.52
1:B:316:ILE:C	1:B:317:ILE:HD12	2.31	0.51
1:B:225:MET:O	1:B:229:VAL:HG23	2.10	0.51
1:B:356:ILE:HG12	1:B:368:PHE:O	2.10	0.51
1:B:282:LYS:HD3	1:B:283:LYS:N	2.26	0.51
1:B:526:PHE:HB3	1:B:553:VAL:HG13	1.93	0.50
2:A:807:MAN:O5	2:A:808:NAG:H82	2.11	0.50
1:B:274:ARG:HH21	1:B:284:HIS:CD2	2.29	0.50
1:A:594:GLY:O	1:A:605:VAL:HG23	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:421:GLU:OE2	1:B:421:GLU:N	2.44	0.49
1:A:256:LYS:HD2	1:A:273:TRP:CZ2	2.47	0.49
1:A:362:SER:O	1:A:365:PRO:HD3	2.12	0.49
1:B:237:ARG:NH2	1:B:422:ASN:HB3	2.28	0.49
1:A:529:ASN:HB3	1:A:548:CYS:SG	2.53	0.49
2:A:805:MAN:O4	1:B:464:SER:HA	2.12	0.49
1:B:287:ALA:HB3	1:B:314:THR:HG22	1.93	0.49
1:A:325:LEU:O	1:A:329:VAL:HG23	2.12	0.48
1:A:335:MET:O	1:A:338:ASN:HB2	2.13	0.48
1:B:465:GLU:HG2	5:B:66:HOH:O	2.12	0.48
1:A:317:ILE:HG23	1:A:330:ILE:CG2	2.43	0.48
1:B:470:GLU:HB3	1:B:480:PRO:HG2	1.95	0.48
1:B:594:GLY:O	1:B:605:VAL:HG23	2.14	0.48
1:A:541:GLU:HG2	1:A:546:ILE:CG1	2.44	0.48
1:A:199:PHE:CD2	1:A:393:PRO:HD2	2.49	0.48
1:A:576:SER:C	1:A:578:GLU:H	2.17	0.48
1:A:230:ASN:O	1:A:234:GLU:HG3	2.14	0.47
1:B:281:ARG:HH11	1:B:281:ARG:HG3	1.79	0.47
1:B:406:ASN:ND2	1:B:408:LEU:HD12	2.29	0.47
1:A:520:GLU:O	1:A:561:LYS:HE3	2.15	0.47
1:A:387:GLU:N	1:A:387:GLU:OE2	2.44	0.47
1:A:541:GLU:HG2	1:A:546:ILE:HG13	1.96	0.47
1:B:303:TYR:CD2	1:B:312:ARG:HA	2.50	0.47
1:B:451:LYS:HD2	1:B:451:LYS:N	2.30	0.46
1:B:341:ILE:HD11	1:B:377:CYS:HB2	1.97	0.46
1:A:588:GLY:C	1:A:596:VAL:HG22	2.35	0.46
1:A:303:TYR:CD2	1:A:312:ARG:HA	2.51	0.46
1:B:604:ASP:OD2	1:B:606:ALA:HB3	2.15	0.46
1:A:567:GLN:OE1	1:A:569:ASN:N	2.49	0.46
1:B:454:LYS:O	1:B:457:THR:HG23	2.16	0.46
1:A:358:ARG:HH11	1:A:358:ARG:HG3	1.79	0.46
1:B:558:LEU:HD12	1:B:559:TYR:N	2.30	0.46
1:A:239:MET:O	1:A:241:ILE:HG13	2.15	0.46
1:A:457:THR:HG21	5:A:146:HOH:O	2.16	0.45
1:A:312:ARG:HD2	1:B:567:GLN:HE22	1.81	0.45
1:A:528:ARG:HG3	1:A:528:ARG:HH11	1.80	0.45
1:B:420:PRO:HG2	1:B:421:GLU:OE2	2.17	0.45
1:A:588:GLY:O	1:A:596:VAL:HG22	2.16	0.45
1:A:356:ILE:HD13	1:A:368:PHE:C	2.37	0.45
1:A:247:GLY:CA	1:A:399:ILE:HD13	2.47	0.45
1:B:274:ARG:HD3	1:B:310:PRO:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:358:ARG:HH21	1:B:358:ARG:HG2	1.83	0.44
2:A:805:MAN:O2	2:A:806:NAG:H82	2.18	0.44
1:A:256:LYS:HD2	1:A:273:TRP:HZ2	1.82	0.44
1:B:414:GLU:OE2	1:B:435:LEU:N	2.48	0.44
1:B:541:GLU:O	1:B:542:ASN:HB2	2.18	0.44
1:A:363:PRO:HA	5:A:91:HOH:O	2.17	0.44
1:A:465:GLU:OE2	2:B:806:NAG:H3	2.17	0.44
1:A:526:PHE:O	1:A:553:VAL:HG13	2.18	0.44
1:B:204:LEU:HD13	1:B:225:MET:HE2	2.00	0.44
1:A:267:LEU:HD22	1:A:301:LEU:HG	1.99	0.44
1:B:405:GLY:O	1:B:418:GLY:HA2	2.17	0.44
1:B:506:MET:HG2	1:B:520:GLU:CD	2.37	0.44
1:A:218:LEU:O	1:A:222:LYS:HG3	2.17	0.44
1:B:202:LEU:HD12	1:B:203:VAL:H	1.83	0.44
1:B:464:SER:HB2	5:B:66:HOH:O	2.17	0.44
1:B:463:MET:HG3	1:B:467:ASP:OD1	2.17	0.44
1:A:356:ILE:HD13	1:A:368:PHE:O	2.18	0.43
1:A:211:VAL:HG21	1:A:252:SER:HA	2.01	0.43
1:B:454:LYS:H	1:B:454:LYS:HE3	1.82	0.43
1:A:450:CYS:O	1:A:451:LYS:HD3	2.18	0.43
1:B:453:SER:O	1:B:474:GLY:HA2	2.19	0.43
1:A:567:GLN:O	1:A:568:ASN:CB	2.66	0.42
1:A:391:ASN:H	1:A:391:ASN:HD22	1.66	0.42
1:B:386:PRO:HB2	1:B:389:ILE:HG12	2.01	0.42
1:B:237:ARG:HE	1:B:407:GLU:CD	2.23	0.42
1:B:330:ILE:O	1:B:333:HIS:HB3	2.19	0.42
1:A:197:PHE:HA	1:A:240:TYR:O	2.19	0.42
1:B:557:ARG:HG2	1:B:573:MET:HG3	2.01	0.42
1:B:259:VAL:CG1	1:B:293:ILE:HD12	2.49	0.42
1:A:529:ASN:ND2	1:A:557:ARG:HB3	2.17	0.42
1:B:281:ARG:NH1	1:B:281:ARG:HG3	2.34	0.42
1:B:235:ILE:HG23	1:B:367:THR:O	2.20	0.42
1:A:574:PHE:CE1	1:A:582:LYS:HG2	2.54	0.42
1:A:317:ILE:HD11	1:A:331:MET:HA	2.02	0.42
1:A:569:ASN:HD22	1:A:570:PRO:HD2	1.85	0.41
1:A:356:ILE:HD12	1:A:356:ILE:N	2.35	0.41
1:A:307:MET:O	1:A:308:CYS:CB	2.67	0.41
1:B:509:GLN:HB3	5:B:117:HOH:O	2.19	0.41
1:B:273:TRP:O	1:B:277:ASP:HB3	2.20	0.41
1:A:385:ASN:O	1:A:387:GLU:N	2.53	0.41
1:B:394:LEU:HB2	1:B:397:ASP:OD2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:487:ASN:ND2	1:B:500:ASN:H	2.18	0.41
1:B:230:ASN:O	1:B:234:GLU:HG3	2.20	0.41
1:B:200:VAL:HG11	1:B:335:MET:CE	2.50	0.41
1:A:529:ASN:ND2	1:A:559:TYR:OH	2.53	0.41
1:A:513:LEU:HD11	1:A:583:GLY:HA2	2.02	0.41
1:A:514:PHE:CZ	1:A:572:LYS:HG3	2.56	0.41
1:B:386:PRO:HB2	1:B:389:ILE:CD1	2.51	0.41
1:B:362:SER:HB2	1:B:363:PRO:HD2	2.03	0.41
1:A:576:SER:C	1:A:578:GLU:N	2.74	0.41
1:A:262:GLU:HB2	5:A:69:HOH:O	2.21	0.41
1:A:414:GLU:HB3	1:A:435:LEU:HG	2.02	0.40
1:A:342:ASN:HB2	1:A:376:GLU:OE1	2.21	0.40
1:B:253:ASN:HB2	5:B:115:HOH:O	2.20	0.40
1:A:386:PRO:O	1:A:389:ILE:HG12	2.21	0.40
1:B:326:VAL:O	1:B:330:ILE:HG13	2.21	0.40
1:A:309:HIS:CE1	1:A:311:LYS:H	2.38	0.40
1:A:275:LYS:HZ3	1:A:310:PRO:HG2	1.87	0.40
1:B:499:TYR:CZ	1:B:500:ASN:ND2	2.89	0.40
1:B:317:ILE:N	1:B:317:ILE:HD12	2.37	0.40
1:B:287:ALA:O	1:B:314:THR:HA	2.21	0.40

There are no symmetry-related clashes.

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	413/419 (99%)	380 (92%)	31 (8%)	2 (0%)	34	63
1	B	413/419 (99%)	381 (92%)	29 (7%)	3 (1%)	26	55
All	All	826/838 (99%)	761 (92%)	60 (7%)	5 (1%)	30	59

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	430	ALA
1	A	568	ASN
1	B	277	ASP
1	B	196	PRO
1	A	566	GLY

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	361/365 (99%)	355 (98%)	6 (2%)	68 90
1	B	361/365 (99%)	347 (96%)	14 (4%)	39 70
All	All	722/730 (99%)	702 (97%)	20 (3%)	51 81

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	262	GLU
1	A	280	THR
1	A	283	LYS
1	A	538	CYS
1	A	569	ASN
1	A	596	VAL
1	B	254	GLU
1	B	360	GLU
1	B	385	ASN
1	B	388	CYS
1	B	391	ASN
1	B	425	ASN
1	B	436	LYS
1	B	451	LYS
1	B	454	LYS
1	B	506	MET
1	B	523	ASP
1	B	538	CYS
1	B	553	VAL

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Mol	Chain	Res	Type
1	B	599	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	288	GLN
1	A	318	GLN
1	A	383	ASN
1	A	422	ASN
1	A	485	HIS
1	A	487	ASN
1	A	500	ASN
1	A	502	ASN
1	A	508	HIS
1	A	529	ASN
1	A	569	ASN
1	A	599	ASN
1	B	268	ASN
1	B	288	GLN
1	B	324	ASN
1	B	385	ASN
1	B	425	ASN
1	B	440	GLN
1	B	475	GLN
1	B	487	ASN
1	B	500	ASN
1	B	529	ASN
1	B	542	ASN
1	B	544	ASN
1	B	563	ASN
1	B	581	HIS
1	B	599	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

18 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	$\#\|Z\|>2$
2	NAG	A	801	1,2	14,14,15	0.55	0	15,19,21	0.83	1 (6%)
2	NAG	A	802	2	14,14,15	0.51	0	15,19,21	0.87	1 (6%)
2	BMA	A	803	2	11,11,12	0.52	0	14,15,17	0.58	0
2	NAG	A	804	2	14,14,15	0.46	0	15,19,21	0.77	1 (6%)
2	MAN	A	805	2	11,11,12	0.92	1 (9%)	14,15,17	1.70	2 (14%)
2	NAG	A	806	2	14,14,15	0.64	0	15,19,21	0.75	1 (6%)
2	MAN	A	807	2	11,11,12	0.64	0	14,15,17	0.63	0
2	NAG	A	808	2	14,14,15	0.56	0	15,19,21	0.70	1 (6%)
2	FUC	A	809	2	10,10,11	0.52	0	14,14,16	0.59	0
2	NAG	B	801	1,2	14,14,15	0.53	0	15,19,21	0.86	1 (6%)
2	NAG	B	802	2	14,14,15	0.54	0	15,19,21	0.76	0
2	BMA	B	803	2	11,11,12	0.52	0	14,15,17	0.63	0
2	NAG	B	804	2	14,14,15	0.44	0	15,19,21	0.82	1 (6%)
2	MAN	B	805	2	11,11,12	0.96	1 (9%)	14,15,17	1.53	1 (7%)
2	NAG	B	806	2	14,14,15	0.67	0	15,19,21	0.78	1 (6%)
2	MAN	B	807	2	11,11,12	0.38	0	14,15,17	0.75	1 (7%)
2	NAG	B	808	2	14,14,15	0.48	0	15,19,21	0.65	0
2	FUC	B	809	2	10,10,11	0.64	0	14,14,16	0.54	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
2	NAG	A	801	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	802	2	-	0/6/23/26	0/1/1/1
2	BMA	A	803	2	-	0/2/19/22	0/1/1/1
2	NAG	A	804	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	A	805	2	-	0/2/19/22	0/1/1/1
2	NAG	A	806	2	-	0/6/23/26	0/1/1/1
2	MAN	A	807	2	-	0/2/19/22	0/1/1/1
2	NAG	A	808	2	-	0/6/23/26	0/1/1/1
2	FUC	A	809	2	-	0/0/17/20	0/1/1/1
2	NAG	B	801	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	802	2	-	0/6/23/26	0/1/1/1
2	BMA	B	803	2	-	0/2/19/22	0/1/1/1
2	NAG	B	804	2	-	0/6/23/26	0/1/1/1
2	MAN	B	805	2	-	0/2/19/22	0/1/1/1
2	NAG	B	806	2	-	0/6/23/26	0/1/1/1
2	MAN	B	807	2	-	0/2/19/22	0/1/1/1
2	NAG	B	808	2	-	0/6/23/26	0/1/1/1
2	FUC	B	809	2	-	0/0/17/20	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	805	MAN	C2-C3	2.45	1.55	1.52
2	A	805	MAN	C2-C3	2.51	1.56	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	804	NAG	C2-N2-C7	-2.57	119.73	123.04
2	A	801	NAG	C2-N2-C7	-2.53	119.79	123.04
2	A	804	NAG	C2-N2-C7	-2.45	119.89	123.04
2	A	806	NAG	C2-N2-C7	-2.20	120.21	123.04
2	B	806	NAG	C2-N2-C7	-2.14	120.30	123.04
2	A	808	NAG	C2-N2-C7	-2.12	120.31	123.04
2	B	801	NAG	C2-N2-C7	-2.08	120.37	123.04
2	A	802	NAG	C2-N2-C7	-2.07	120.38	123.04
2	B	807	MAN	C1-O5-C5	2.04	114.83	112.25
2	A	805	MAN	C2-C3-C4	2.14	114.68	111.04
2	B	805	MAN	C1-C2-C3	4.77	115.18	109.54
2	A	805	MAN	C1-C2-C3	5.53	116.08	109.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	804	NAG	1	0
2	A	805	MAN	2	0
2	A	806	NAG	3	0
2	A	807	MAN	1	0
2	A	808	NAG	1	0
2	B	806	NAG	1	0

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

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 (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, 95^{th} 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	415/419 (99%)	-0.17	3 (0%)	89	90	6, 22, 43, 66	0
1	B	415/419 (99%)	-0.12	3 (0%)	89	90	8, 24, 46, 61	0
All	All	830/838 (99%)	-0.14	6 (0%)	89	90	6, 23, 44, 66	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	568	ASN	3.8
1	A	567	GLN	3.4
1	B	542	ASN	3.0
1	B	424	GLN	2.8
1	A	352	ASP	2.6
1	B	568	ASN	2.5

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

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, 95^{th} 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
2	NAG	A	806	14/15	0.91	0.19	1.29	14,31,33,34	0
2	NAG	B	804	14/15	0.93	0.16	0.49	24,31,34,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	NAG	B	801	14/15	0.92	0.17	0.21	29,34,41,46	0
2	NAG	B	806	14/15	0.89	0.17	0.07	28,35,38,39	0
2	NAG	A	804	14/15	0.94	0.16	0.06	30,35,40,40	0
2	NAG	A	801	14/15	0.94	0.15	-0.01	26,31,35,37	0
2	FUC	A	809	10/11	0.93	0.23	-	23,30,34,39	0
2	NAG	A	802	14/15	0.93	0.14	-	24,32,35,37	0
2	FUC	B	809	10/11	0.92	0.21	-	26,30,34,36	0
2	BMA	A	803	11/12	0.94	0.15	-	29,35,38,39	0
2	MAN	A	807	11/12	0.87	0.17	-	32,38,45,46	0
2	NAG	B	808	14/15	0.86	0.26	-	42,47,48,51	0
2	NAG	B	802	14/15	0.93	0.18	-	24,39,45,46	0
2	BMA	B	803	11/12	0.96	0.20	-	34,40,42,45	0
2	MAN	A	805	11/12	0.90	0.24	-	34,37,39,48	0
2	MAN	B	805	11/12	0.89	0.24	-	32,40,43,45	0
2	MAN	B	807	11/12	0.93	0.18	-	33,36,39,40	0
2	NAG	A	808	14/15	0.91	0.30	-	42,45,48,49	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, 95^{th} 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	CA	B	711	1/1	0.96	0.12	-0.50	23,23,23,23	0
4	CA	A	703	1/1	0.98	0.13	-1.64	23,23,23,23	0
4	CA	B	713	1/1	0.98	0.12	-1.70	20,20,20,20	0
4	CA	A	702	1/1	0.97	0.06	-2.77	20,20,20,20	0
4	CA	A	701	1/1	0.99	0.06	-2.84	21,21,21,21	0
4	CA	B	712	1/1	0.99	0.06	-3.14	22,22,22,22	0
3	ZN	A	700	1/1	1.00	0.10	-3.56	18,18,18,18	0
3	ZN	B	700	1/1	1.00	0.12	-	17,17,17,17	0

6.5 Other polymers (i)

There are no such residues in this entry.