Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:33 AM GMT

PDB ID	:	3F6K
Title	:	Crystal structure of the Vps10p domain of human sortilin/NTS3 in complex with neurotensin
Authors	:	Quistgaard, E.M.; Madsen, P.; Groftehauge, M.K.; Nissen, P.; Petersen, C.M.; Thirup, S.
Deposited on	:	2008-11-06
Resolution	:	2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

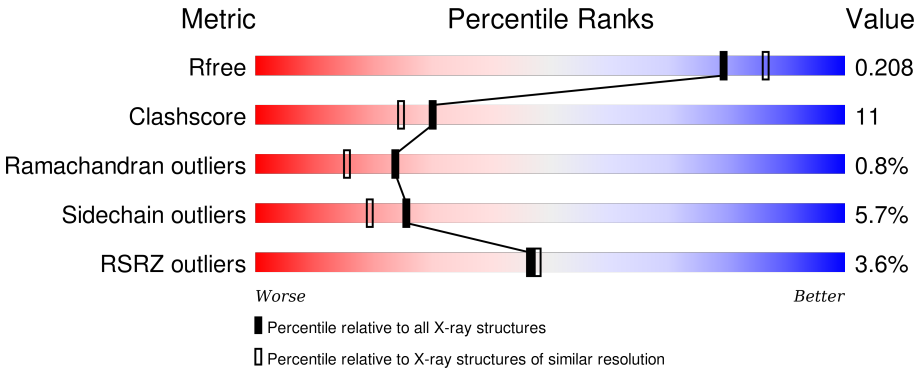
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	685	4%, 73%, 19%
2	N	13	23%, 8%, 69%

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
5	PGE	A	951	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5600 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 Sortilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	657	Total	C	N	O	S	0	0	0
			5176	3268	871	1009	28			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	617	MET	VAL	VARIANT	UNP Q99523
A	724	HIS	-	EXPRESSION TAG	UNP Q99523
A	725	HIS	-	EXPRESSION TAG	UNP Q99523
A	726	HIS	-	EXPRESSION TAG	UNP Q99523
A	727	HIS	-	EXPRESSION TAG	UNP Q99523
A	728	HIS	-	EXPRESSION TAG	UNP Q99523
A	729	HIS	-	EXPRESSION TAG	UNP Q99523

- Molecule 2 is a protein called Neurotensin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	N	4	Total	C	N	O	0	0	0
			36	26	4	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
N	1	PCA	GLN	SEE REMARK 999	UNP P30990

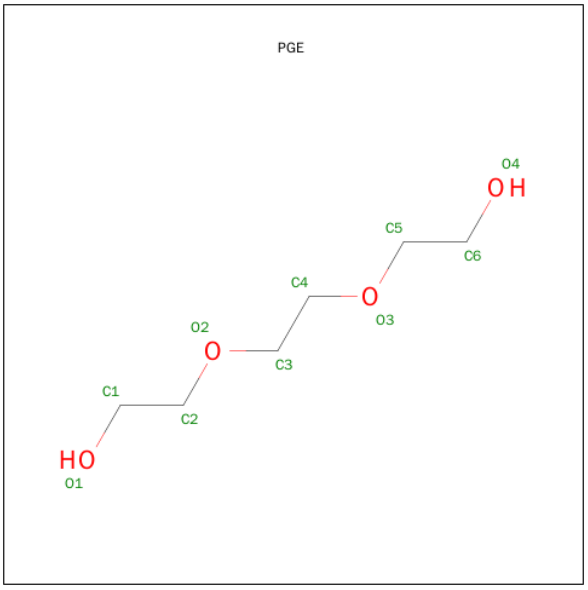
- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

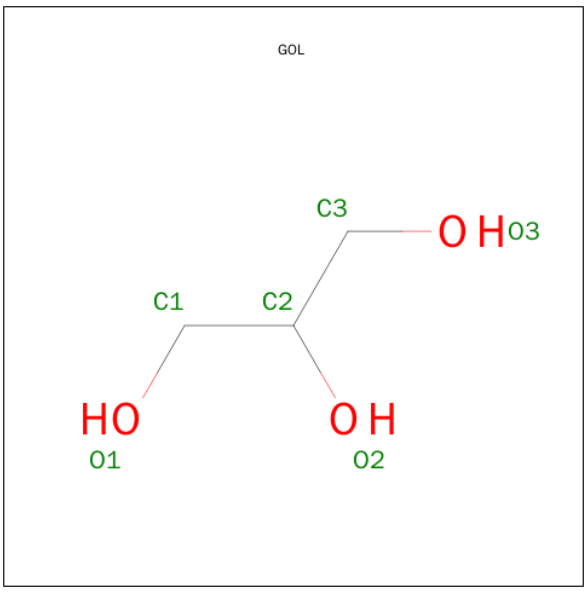
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 7 is water.

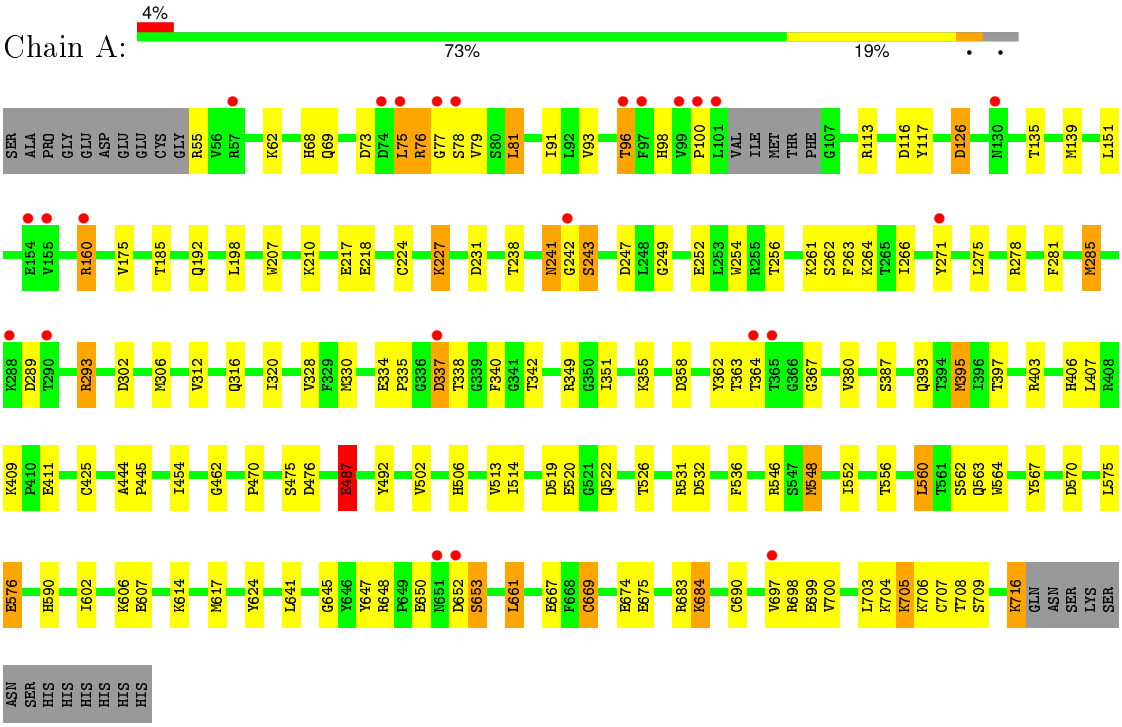
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	304	Total O 304 304	0	0
7	N	1	Total O 1 1	0	0

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: Sortilin

Chain A:



- Molecule 2: Neurotensin

Chain N:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	145.76Å 74.53Å 108.33Å 90.00° 131.87° 90.00°	Depositor
Resolution (Å)	40.32 – 2.00 39.52 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.4 (40.32-2.00) 97.6 (39.52-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.43 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.204 , 0.229 0.204 , 0.208	Depositor DCC
R_{free} test set	2876 reflections (5.31%)	DCC
Wilson B-factor (Å^2)	29.3	Xtriage
Anisotropy	0.443	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 60.0	EDS
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 66579 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5600	wwPDB-VP
Average B, all atoms (Å^2)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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: GOL, BMA, PGE, 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	0.98	6/5298 (0.1%)	0.91	11/7176 (0.2%)
2	N	0.78	0/37	1.30	0/48
All	All	0.98	6/5335 (0.1%)	0.92	11/7224 (0.2%)

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
1	A	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	487	GLU	CG-CD	7.31	1.62	1.51
1	A	224	CYS	CB-SG	-6.84	1.70	1.82
1	A	698	ARG	CZ-NH1	6.02	1.40	1.33
1	A	487	GLU	CB-CG	5.90	1.63	1.52
1	A	492	TYR	CD1-CE1	5.39	1.47	1.39
1	A	475	SER	CA-CB	5.10	1.60	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	293	ARG	CB-CA-C	-6.99	96.43	110.40
1	A	570	ASP	CB-CG-OD1	6.71	124.34	118.30
1	A	293	ARG	NE-CZ-NH2	-6.56	117.02	120.30
1	A	698	ARG	NE-CZ-NH2	-6.34	117.13	120.30
1	A	570	ASP	CB-CG-OD2	-5.94	112.96	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	113	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	A	395	MET	CG-SD-CE	-5.81	90.90	100.20
1	A	81	LEU	CA-CB-CG	5.32	127.54	115.30
1	A	548	MET	CG-SD-CE	-5.24	91.81	100.20
1	A	126	ASP	CB-CG-OD1	5.22	122.99	118.30
1	A	606	LYS	CB-CA-C	-5.04	100.31	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	652	ASP	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5176	0	4965	114	0
2	N	36	0	38	2	0
3	A	28	0	25	0	0
4	A	39	0	34	0	0
5	A	10	0	14	1	0
6	A	6	0	8	2	0
7	A	304	0	0	11	0
7	N	1	0	0	0	0
All	All	5600	0	5084	114	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 11.

All (114) 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:648:ARG:HD2	1:A:653:SER:HA	1.30	1.06
1:A:96:THR:HG21	1:A:135:THR:OG1	1.58	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:548:MET:HE3	1:A:575:LEU:HD12	1.44	0.98
1:A:502:VAL:HG11	1:A:514:ILE:HD11	1.44	0.96
1:A:502:VAL:CG1	1:A:514:ILE:HD11	1.98	0.93
1:A:293:ARG:HH22	1:A:667:GLU:CD	1.82	0.82
1:A:661:LEU:CD2	1:A:707:CYS:SG	2.67	0.81
1:A:81:LEU:HD11	1:A:91:ILE:HG21	1.62	0.80
1:A:81:LEU:HD13	1:A:93:VAL:HG22	1.64	0.78
1:A:519:ASP:OD2	1:A:522:GLN:HB2	1.84	0.77
1:A:648:ARG:HD2	1:A:653:SER:CA	2.12	0.76
1:A:75:LEU:HD23	1:A:79:VAL:HG21	1.69	0.74
1:A:403:ARG:HH11	1:A:403:ARG:HG3	1.53	0.74
1:A:96:THR:CG2	1:A:135:THR:OG1	2.37	0.73
1:A:560:LEU:H	1:A:560:LEU:HD12	1.54	0.72
1:A:81:LEU:HD11	1:A:91:ILE:CG2	2.21	0.71
1:A:487:GLU:CD	1:A:487:GLU:H	1.95	0.70
1:A:648:ARG:CD	1:A:653:SER:HA	2.18	0.69
1:A:320:ILE:HD13	1:A:330:MET:SD	2.33	0.69
1:A:669:CYS:HG	1:A:707:CYS:HG	0.74	0.68
1:A:513:VAL:CG1	1:A:526:THR:HG23	2.23	0.68
1:A:81:LEU:CD1	1:A:91:ILE:CG2	2.72	0.68
1:A:513:VAL:HG13	1:A:526:THR:HG23	1.75	0.67
1:A:661:LEU:HD22	1:A:707:CYS:SG	2.34	0.66
1:A:506:HIS:CE1	5:A:951:PGE:H1	2.32	0.65
1:A:334:GLU:HG3	1:A:335:PRO:HD2	1.79	0.65
1:A:546:ARG:NH1	1:A:624:TYR:O	2.23	0.65
1:A:684:LYS:HE2	1:A:690:CYS:SG	2.38	0.64
1:A:409:LYS:HD3	1:A:425:CYS:SG	2.37	0.64
1:A:192:GLN:OE1	1:A:231:ASP:HA	1.99	0.62
1:A:75:LEU:HD21	1:A:79:VAL:HG11	1.81	0.62
1:A:699:GLU:HA	1:A:699:GLU:OE1	1.99	0.62
1:A:238:THR:OG1	1:A:252:GLU:HG3	1.99	0.61
1:A:395:MET:HA	1:A:395:MET:HE2	1.84	0.60
1:A:241:ASN:HB2	1:A:247:ASP:OD2	2.01	0.60
1:A:444:ALA:HB1	1:A:445:PRO:CD	2.33	0.59
1:A:81:LEU:CD1	1:A:93:VAL:HG22	2.33	0.58
1:A:312:VAL:HB	1:A:316:GLN:HB2	1.87	0.57
1:A:355:LYS:HD3	1:A:358:ASP:OD2	2.04	0.57
1:A:160:ARG:HG3	1:A:160:ARG:HH11	1.70	0.56
1:A:520:GLU:OE1	1:A:614:LYS:HE2	2.04	0.56
1:A:454:ILE:HD13	1:A:476:ASP:HA	1.86	0.56
1:A:363:THR:O	1:A:367:GLY:HA2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:THR:HG22	7:A:1275:HOH:O	2.05	0.55
1:A:661:LEU:HD23	1:A:707:CYS:SG	2.47	0.55
1:A:607:GLU:OE1	7:A:1270:HOH:O	2.18	0.54
1:A:78:SER:HB2	1:A:96:THR:HG23	1.90	0.54
1:A:337:ASP:HA	7:A:1269:HOH:O	2.08	0.54
1:A:548:MET:SD	7:A:1277:HOH:O	2.59	0.54
1:A:293:ARG:NH2	1:A:667:GLU:CD	2.57	0.54
1:A:645:GLY:HA3	1:A:703:LEU:HD13	1.89	0.53
1:A:556:THR:O	1:A:562:SER:HB2	2.08	0.53
1:A:576:GLU:OE1	6:A:952:GOL:C1	2.56	0.53
1:A:77:GLY:HA3	1:A:96:THR:O	2.09	0.52
1:A:185:THR:HB	7:A:1240:HOH:O	2.09	0.52
1:A:387:SER:HB2	1:A:393:GLN:OE1	2.08	0.52
1:A:706:LYS:HG2	1:A:707:CYS:SG	2.50	0.51
1:A:395:MET:HE1	1:A:406:HIS:ND1	2.25	0.51
1:A:81:LEU:HD12	1:A:91:ILE:CG2	2.41	0.51
1:A:264:LYS:HD3	1:A:266:ILE:CG2	2.41	0.50
1:A:548:MET:CE	1:A:575:LEU:HD12	2.31	0.49
1:A:705:LYS:O	1:A:709:SER:HB2	2.13	0.49
1:A:75:LEU:HB3	1:A:564:TRP:CD1	2.47	0.49
1:A:520:GLU:OE1	1:A:614:LYS:CE	2.61	0.49
1:A:320:ILE:CD1	1:A:330:MET:CE	2.91	0.49
1:A:337:ASP:HA	7:A:1237:HOH:O	2.12	0.49
1:A:207:TRP:CE2	1:A:218:GLU:HB3	2.48	0.48
1:A:380:VAL:HA	1:A:397:THR:O	2.14	0.47
1:A:716:LYS:HE2	1:A:716:LYS:HB3	1.58	0.47
1:A:502:VAL:CG1	1:A:514:ILE:CD1	2.84	0.47
1:A:242:GLY:O	1:A:243:SER:O	2.33	0.47
1:A:617:MET:HE1	7:A:1287:HOH:O	2.15	0.47
1:A:704:LYS:HE3	1:A:708:THR:HG21	1.97	0.47
1:A:271:TYR:CD1	1:A:285:MET:SD	3.08	0.47
1:A:227:LYS:NZ	2:N:11:TYR:HE1	2.14	0.46
1:A:502:VAL:HG13	1:A:514:ILE:HD11	1.92	0.46
1:A:264:LYS:HE3	1:A:302:ASP:OD1	2.16	0.46
1:A:281:PHE:HZ	1:A:328:VAL:HG11	1.81	0.46
1:A:641:LEU:HG	1:A:683:ARG:O	2.16	0.45
1:A:198:LEU:C	1:A:198:LEU:HD12	2.36	0.45
1:A:338:THR:HB	1:A:340:PHE:CE2	2.52	0.45
1:A:160:ARG:HG3	1:A:160:ARG:NH1	2.32	0.45
1:A:210:LYS:NZ	1:A:217:GLU:OE2	2.50	0.45
1:A:513:VAL:HG11	1:A:526:THR:HG23	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:411:GLU:HG2	7:A:1133:HOH:O	2.17	0.44
1:A:139:MET:HG2	1:A:151:LEU:CD2	2.47	0.44
1:A:227:LYS:NZ	2:N:11:TYR:CE1	2.86	0.44
1:A:395:MET:CA	1:A:395:MET:HE2	2.46	0.43
1:A:556:THR:CG2	1:A:563:GLN:HG2	2.48	0.43
1:A:81:LEU:HD13	1:A:93:VAL:CG2	2.42	0.43
1:A:249:GLY:HA2	1:A:285:MET:HG3	2.00	0.43
1:A:62:LYS:NZ	7:A:1120:HOH:O	2.44	0.43
1:A:266:ILE:HA	7:A:1261:HOH:O	2.18	0.43
1:A:264:LYS:HD3	1:A:266:ILE:HG22	2.00	0.43
1:A:271:TYR:HD1	1:A:285:MET:SD	2.41	0.43
1:A:425:CYS:HB2	1:A:462:GLY:O	2.19	0.42
1:A:116:ASP:O	1:A:117:TYR:HB2	2.19	0.42
1:A:256:THR:O	1:A:256:THR:HG23	2.19	0.42
1:A:590:HIS:CG	1:A:602:ILE:HG13	2.55	0.42
1:A:69:GLN:HG3	1:A:567:TYR:CE1	2.55	0.42
1:A:444:ALA:HB1	1:A:445:PRO:HD2	2.01	0.42
1:A:560:LEU:CD1	1:A:560:LEU:H	2.26	0.42
1:A:76:ARG:HD3	1:A:76:ARG:HA	1.62	0.42
1:A:407:LEU:N	1:A:407:LEU:HD12	2.34	0.42
1:A:395:MET:CE	1:A:406:HIS:ND1	2.83	0.42
1:A:403:ARG:HG3	1:A:403:ARG:NH1	2.28	0.41
1:A:674:GLU:HG2	1:A:675:GLU:N	2.35	0.41
1:A:576:GLU:OE1	6:A:952:GOL:H11	2.19	0.41
1:A:98:HIS:O	1:A:100:PRO:HD3	2.21	0.41
1:A:68:HIS:HE1	7:A:1135:HOH:O	2.04	0.41
1:A:556:THR:HG22	1:A:563:GLN:HG2	2.02	0.41
1:A:254:TRP:CE3	1:A:263:PHE:HB3	2.56	0.41
1:A:536:PHE:HE1	1:A:552:ILE:HG23	1.87	0.40
1:A:349:ARG:HB3	1:A:351:ILE:HD12	2.02	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	653/685 (95%)	621 (95%)	27 (4%)	5 (1%)	24 15
2	N	2/13 (15%)	2 (100%)	0	0	100 100
All	All	655/698 (94%)	623 (95%)	27 (4%)	5 (1%)	24 15

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	241	ASN
1	A	243	SER
1	A	653	SER
1	A	650	GLU
1	A	337	ASP

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	574/599 (96%)	542 (94%)	32 (6%)	26 20
2	N	4/12 (33%)	3 (75%)	1 (25%)	1 0
All	All	578/611 (95%)	545 (94%)	33 (6%)	25 19

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

Mol	Chain	Res	Type
1	A	55	ARG
1	A	73	ASP
1	A	75	LEU
1	A	76	ARG
1	A	96	THR
1	A	126	ASP
1	A	160	ARG
1	A	175	VAL

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Mol	Chain	Res	Type
1	A	227	LYS
1	A	261	LYS
1	A	262	SER
1	A	275	LEU
1	A	278	ARG
1	A	285	MET
1	A	289	ASP
1	A	306	MET
1	A	362	TYR
1	A	364	THR
1	A	470	PRO
1	A	487	GLU
1	A	531	ARG
1	A	532	ASP
1	A	560	LEU
1	A	576	GLU
1	A	647	TYR
1	A	661	LEU
1	A	669	CYS
1	A	684	LYS
1	A	697	VAL
1	A	700	VAL
1	A	705	LYS
1	A	716	LYS
2	N	11	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	390	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)

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	801	1,3	14,14,15	1.34	3 (21%)	15,19,21	1.78	3 (20%)
3	NAG	A	802	3	14,14,15	0.69	0	15,19,21	1.17	2 (13%)
4	NAG	A	901	1,4	14,14,15	1.13	2 (14%)	15,19,21	1.61	3 (20%)
4	NAG	A	902	4	14,14,15	0.95	1 (7%)	15,19,21	1.31	1 (6%)
4	BMA	A	903	4	11,11,12	0.67	0	14,15,17	1.09	1 (7%)

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	801	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	802	3	-	0/6/23/26	0/1/1/1
4	NAG	A	901	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	902	4	-	0/6/23/26	0/1/1/1
4	BMA	A	903	4	-	0/2/19/22	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	NAG	C2-N2	-2.62	1.41	1.46
3	A	801	NAG	C3-C2	2.19	1.57	1.52
4	A	902	NAG	C1-C2	2.19	1.55	1.52
3	A	801	NAG	C2-N2	2.33	1.50	1.46
4	A	901	NAG	C1-C2	2.46	1.55	1.52
3	A	801	NAG	C1-C2	3.02	1.56	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	801	NAG	C2-N2-C7	-4.39	117.40	123.04
4	A	901	NAG	C3-C2-N2	-3.40	102.42	110.56
4	A	901	NAG	C6-C5-C4	-2.26	107.45	113.02
3	A	802	NAG	C2-N2-C7	-2.17	120.25	123.04
3	A	801	NAG	C3-C4-C5	-2.04	106.65	110.20
4	A	903	BMA	C1-O5-C5	2.45	115.36	112.25
3	A	802	NAG	C1-O5-C5	2.54	115.47	112.25
3	A	801	NAG	C4-C3-C2	2.88	115.71	111.23
4	A	901	NAG	C1-O5-C5	3.36	116.51	112.25
4	A	902	NAG	C4-C3-C2	3.48	116.64	111.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry (i)

2 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	$\#\|Z\| > 2$
5	PGE	A	951	-	9,9,9	1.04	0	8,8,8	0.27	0
6	GOL	A	952	-	5,5,5	0.22	0	5,5,5	0.64	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	PGE	A	951	-	-	0/7/7/7	0/0/0/0
6	GOL	A	952	-	-	0/4/4/4	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.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	951	PGE	1	0
6	A	952	GOL	2	0

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, 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	657/685 (95%)	0.06	24 (3%) 45 47	12, 28, 46, 58	0
2	N	4/13 (30%)	0.63	0 100 100	52, 53, 61, 62	0
All	All	661/698 (94%)	0.06	24 (3%) 46 48	12, 28, 47, 62	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	99	VAL	8.4
1	A	97	PHE	5.0
1	A	100	PRO	4.9
1	A	101	LEU	4.1
1	A	242	GLY	3.7
1	A	365	THR	3.7
1	A	697	VAL	3.6
1	A	75	LEU	3.6
1	A	77	GLY	3.4
1	A	652	ASP	3.3
1	A	57	ARG	3.2
1	A	96	THR	2.7
1	A	154	GLU	2.7
1	A	651	ASN	2.7
1	A	78	SER	2.6
1	A	155	VAL	2.5
1	A	337	ASP	2.5
1	A	364	THR	2.4
1	A	130	ASN	2.3
1	A	74	ASP	2.3
1	A	271	TYR	2.2
1	A	160	ARG	2.1
1	A	290	THR	2.1
1	A	288	LYS	2.0

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, 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
3	NAG	A	801	14/15	0.96	0.13	-0.29	24,28,30,32	0
4	NAG	A	901	14/15	0.96	0.07	-1.99	25,30,38,46	0
4	BMA	A	903	11/12	0.75	0.33	-	79,83,84,86	0
4	NAG	A	902	14/15	0.87	0.17	-	52,58,64,72	0
3	NAG	A	802	14/15	0.87	0.12	-	41,47,59,60	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
5	PGE	A	951	10/10	0.84	0.16	4.17	66,69,70,70	0
6	GOL	A	952	6/6	0.70	0.20	0.83	60,64,65,66	0

6.5 Other polymers (i)

There are no such residues in this entry.