



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2023 – 06:49 PM JST

PDB ID : 7VF3
Title : Plexin B1 extracellular fragment in complex with lasso-grafted PB1m7 peptide
Authors : Sugano, N.N.; Hirata, K.; Yamashita, K.; Yamamoto, M.; Arimori, T.; Takagi, J.
Deposited on : 2021-09-10
Resolution : 2.29 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

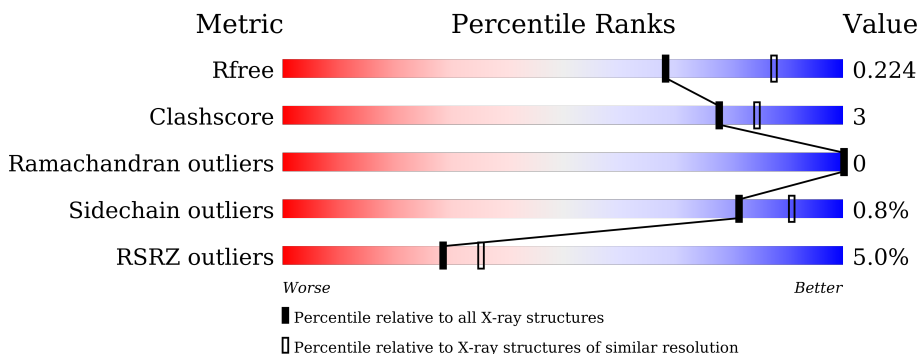
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	525	 4% 87% 7% 6%
1	C	525	 5% 87% 7% 6%
2	B	155	 10% 90%
2	D	155	 2% 9% 90%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7997 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 Plexin-B1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	494	3751	2357	656	716	22	0	0	
1	C	494	3745	2354	653	716	3	0	0	

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	SER	THR	engineered mutation	UNP O43157
A	536	THR	-	expression tag	UNP O43157
A	537	ARG	-	expression tag	UNP O43157
A	538	GLU	-	expression tag	UNP O43157
A	539	ASN	-	expression tag	UNP O43157
A	540	LEU	-	expression tag	UNP O43157
A	541	TYR	-	expression tag	UNP O43157
A	542	PHE	-	expression tag	UNP O43157
A	543	GLN	-	expression tag	UNP O43157
C	19	SER	THR	engineered mutation	UNP O43157
C	536	THR	-	expression tag	UNP O43157
C	537	ARG	-	expression tag	UNP O43157
C	538	GLU	-	expression tag	UNP O43157
C	539	ASN	-	expression tag	UNP O43157
C	540	LEU	-	expression tag	UNP O43157
C	541	TYR	-	expression tag	UNP O43157
C	542	PHE	-	expression tag	UNP O43157
C	543	GLN	-	expression tag	UNP O43157

- Molecule 2 is a protein called Uteroglobin,PB1m7 peptide,Uteroglobin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	15	129	83	20	24	2	0	0	0

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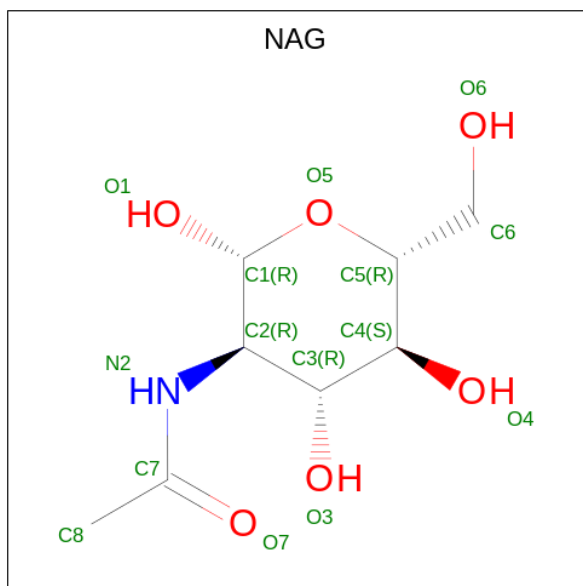
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	15	129	83	20	24	2	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	29	SER	ALA	engineered mutation	UNP P11684
D	29	SER	ALA	engineered mutation	UNP P11684

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



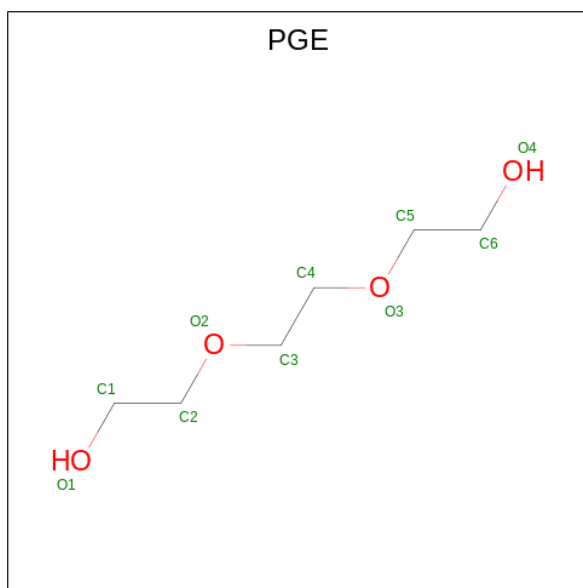
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	14	8	1	5	0	0
3	C	1	14	8	1	5	0	0

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



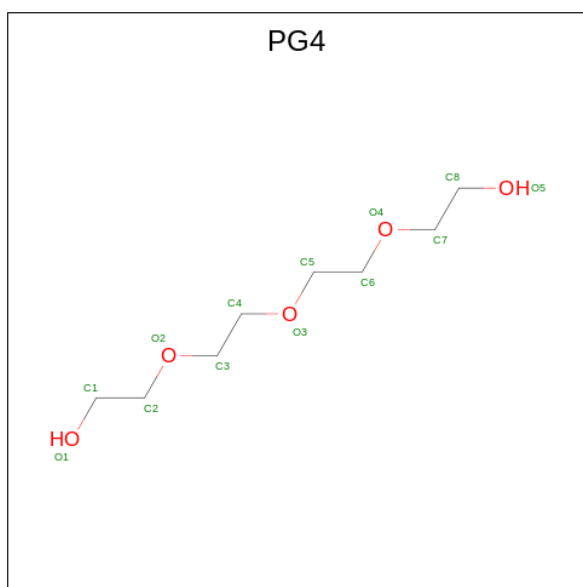
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		

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



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

- Molecule 6 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



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

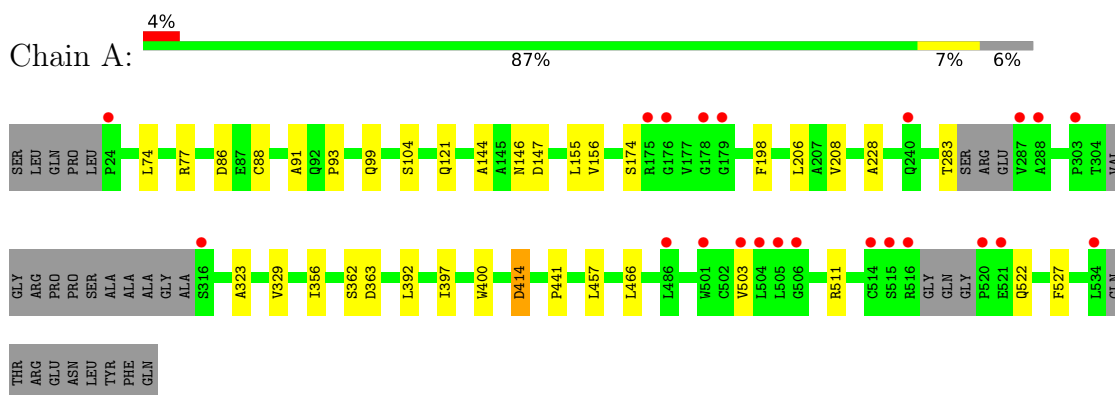
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	64	Total O 64 64	0	0
7	B	2	Total O 2 2	0	0
7	C	77	Total O 77 77	0	0
7	D	3	Total O 3 3	0	0

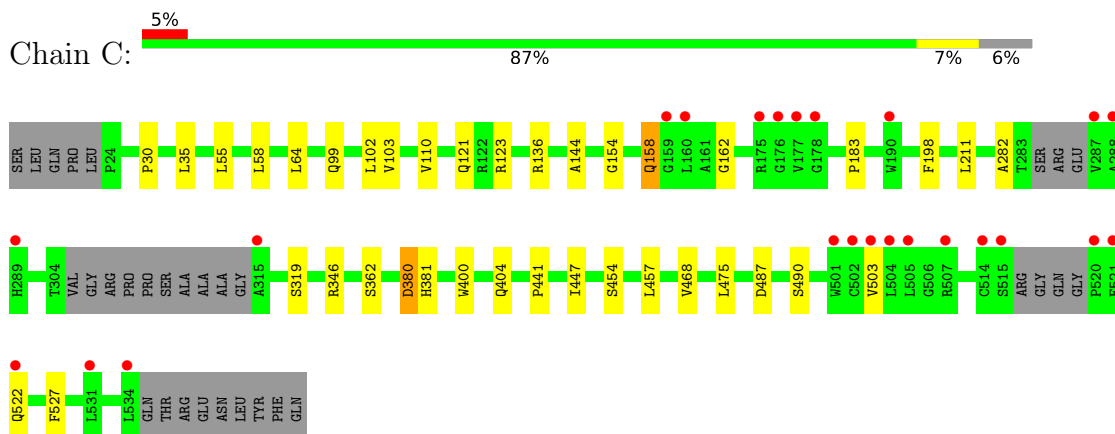
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

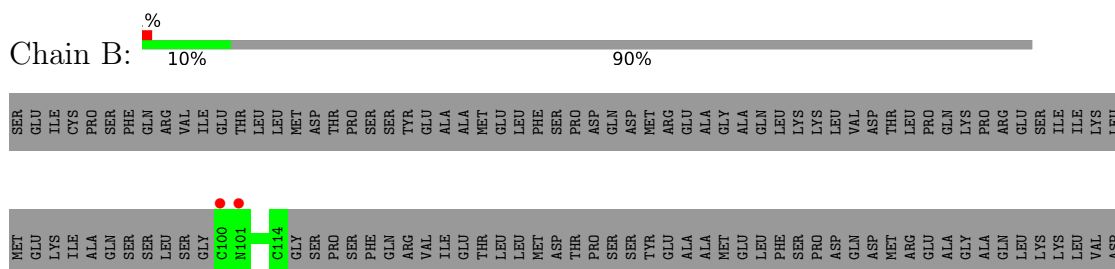
- Molecule 1: Plexin-B1



- Molecule 1: Plexin-B1



- Molecule 2: Uteroglobin, PB1m7 peptide, Uteroglobin



THR
LEU
PRO
GLN
LYS
PRO
ARG
GLU
SER
ILE
ILE
LYS
LEU
MET
GLU
LYS
ILE
ALA
GLN
SER
SER
LEU
CYS
ASN

● Molecule 2: Uteroglobin,PB1m7 peptide,Uteroglobin

Chain D:  2%
9% . 90%

SER
GLU
ILE
CYS
PRO
SER
PHE
GLN
ARG
VAL
ILE
GLU
THR
LEU
LEU
MET
ASP
THR
PRO
SER
TYR
GLU
ALA
ALA
MET
GLU
PHE
SER
PRO
ASP
GLN
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C100
M101
S102
M112
Y113
C114
GLY
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ASP
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GLN
LYS
PRO
ARG
GLU
SER
ILE
ILE
LYS
LEU
MET
GLU
LYS
ILE
ALA
GLN
SER
SER
LEU
CYS
ASN

4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	263.59Å 263.59Å 108.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.41 – 2.29 47.41 – 2.29	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.41-2.29) 99.7 (47.41-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.196 , 0.226 0.193 , 0.224	Depositor DCC
R_{free} test set	4229 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	44.9	Xtrriage
Anisotropy	0.446	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7997	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, PEG, NAG, PG4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.39	0/3846	0.62	0/5252
1	C	0.40	0/3840	0.64	2/5245 (0.0%)
2	B	0.35	0/134	0.45	0/184
2	D	0.39	0/134	0.44	0/184
All	All	0.39	0/7954	0.62	2/10865 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	380	ASP	CB-CG-OD1	6.63	124.27	118.30
1	C	380	ASP	CB-CG-OD2	-5.70	113.17	118.30

There are no chirality outliers.

There are no planarity outliers.

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	3751	0	3601	21	0
1	C	3745	0	3593	23	0
2	B	129	0	107	0	0
2	D	129	0	107	1	0
3	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	14	0	13	0	0
4	A	7	0	10	0	0
5	C	10	0	14	1	0
6	C	52	0	72	5	0
7	A	64	0	0	1	0
7	B	2	0	0	0	0
7	C	77	0	0	0	0
7	D	3	0	0	0	0
All	All	7997	0	7530	45	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 3.

All (45) 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:C:503:VAL:HG21	1:C:522:GLN:HE21	1.40	0.86
1:A:77:ARG:NH1	1:A:146:ASN:O	2.19	0.76
1:C:183:PRO:HD3	1:C:211:LEU:HD11	1.67	0.74
1:C:346:ARG:HE	5:C:602:PGE:H32	1.53	0.72
1:A:363:ASP:OD2	1:C:123:ARG:NH2	2.25	0.70
1:A:77:ARG:HH12	1:A:147:ASP:HA	1.57	0.68
1:C:503:VAL:HG21	1:C:522:GLN:NE2	2.09	0.64
1:C:58:LEU:HD23	1:C:64:LEU:HA	1.78	0.64
1:A:414:ASP:OD2	1:A:511:ARG:NH1	2.32	0.62
1:A:457:LEU:HG	1:A:466:LEU:HD11	1.85	0.59
1:C:487:ASP:OD1	1:C:490:SER:HB2	2.04	0.57
1:C:447:ILE:HD12	1:C:475:LEU:HD12	1.86	0.57
1:A:323:ALA:HB2	1:A:397:ILE:HD11	1.89	0.54
1:A:74:LEU:HD23	1:A:93:PRO:HD3	1.91	0.53
1:C:55:LEU:HD11	1:C:110:VAL:HG11	1.90	0.53
1:C:362:SER:HB2	6:C:605:PG4:H82	1.91	0.53
1:C:103:VAL:HG12	6:C:603:PG4:H72	1.91	0.53
1:A:414:ASP:OD2	1:A:511:ARG:HD2	2.10	0.52
1:C:400:TRP:CZ2	1:C:441:PRO:HD3	2.45	0.52
1:A:99:GLN:HG2	1:A:144:ALA:HB1	1.93	0.50
1:A:206:LEU:HG	1:A:208:VAL:HG23	1.93	0.50
1:A:356:ILE:HG21	1:A:362:SER:HB3	1.93	0.49
1:A:86:ASP:OD2	1:A:86:ASP:N	2.43	0.49
1:A:400:TRP:CZ2	1:A:441:PRO:HD3	2.47	0.49
1:A:329:VAL:HG22	1:A:392:LEU:HD21	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:457:LEU:CD1	1:C:468:VAL:HG13	2.43	0.48
1:C:158:GLN:HB2	1:C:162:GLY:HA2	1.95	0.47
1:A:77:ARG:HH12	1:A:147:ASP:CA	2.28	0.46
1:C:282:ALA:HB1	6:C:604:PG4:H22	1.98	0.45
1:C:102:LEU:HD11	1:C:154:GLY:HA3	1.99	0.45
1:C:319:SER:HG	1:C:404:GLN:HA	1.81	0.45
1:A:121:GLN:HG3	1:A:198:PHE:HZ	1.82	0.45
1:A:88:CYS:SG	1:A:91:ALA:HB2	2.57	0.44
1:A:503:VAL:HG23	1:A:522:GLN:O	2.17	0.44
1:C:30:PRO:HG3	1:C:35:LEU:HD21	2.00	0.44
1:C:121:GLN:HG3	1:C:198:PHE:HZ	1.84	0.43
1:A:155:LEU:HD21	1:A:228:ALA:HB1	2.00	0.43
1:C:99:GLN:HG2	1:C:144:ALA:HB1	2.00	0.43
1:C:487:ASP:OD1	1:C:487:ASP:N	2.51	0.43
1:A:104:SER:HB2	1:A:156:VAL:HG11	2.01	0.43
6:C:604:PG4:H51	6:C:604:PG4:H72	1.73	0.43
6:C:603:PG4:H51	6:C:603:PG4:H31	1.87	0.42
1:A:99:GLN:NE2	7:A:705:HOH:O	2.53	0.41
1:C:380:ASP:OD1	1:C:381:HIS:CE1	2.74	0.41
1:C:454:SER:HB2	2:D:112:TRP:CH2	2.57	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	486/525 (93%)	476 (98%)	10 (2%)	0	100	100
1	C	486/525 (93%)	478 (98%)	8 (2%)	0	100	100
2	B	13/155 (8%)	13 (100%)	0	0	100	100
2	D	13/155 (8%)	12 (92%)	1 (8%)	0	100	100
All	All	998/1360 (73%)	979 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	401/424 (95%)	397 (99%)	4 (1%)	76	87
1	C	400/424 (94%)	397 (99%)	3 (1%)	81	91
2	B	15/141 (11%)	15 (100%)	0	100	100
2	D	15/141 (11%)	15 (100%)	0	100	100
All	All	831/1130 (74%)	824 (99%)	7 (1%)	81	91

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	174	SER
1	A	283	THR
1	A	414	ASP
1	A	527	PHE
1	C	136	ARG
1	C	158	GLN
1	C	527	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	GLN
1	C	256	GLN
1	C	522	GLN

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](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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
6	PG4	C	603	-	12,12,12	0.24	0	11,11,11	0.56	0
5	PGE	C	602	-	9,9,9	0.39	0	8,8,8	0.34	0
6	PG4	C	604	-	12,12,12	0.19	0	11,11,11	0.60	0
4	PEG	A	602	-	6,6,6	0.24	0	5,5,5	0.07	0
6	PG4	C	606	-	12,12,12	0.26	0	11,11,11	0.51	0
3	NAG	A	601	1	14,14,15	0.74	1 (7%)	17,19,21	0.66	1 (5%)
6	PG4	C	605	-	12,12,12	0.26	0	11,11,11	0.49	0
3	NAG	C	601	1	14,14,15	0.48	0	17,19,21	0.50	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PG4	C	603	-	-	1/10/10/10	-
5	PGE	C	602	-	-	4/7/7/7	-
6	PG4	C	604	-	-	7/10/10/10	-
4	PEG	A	602	-	-	2/4/4/4	-
6	PG4	C	606	-	-	5/10/10/10	-
3	NAG	A	601	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PG4	C	605	-	-	5/10/10/10	-
3	NAG	C	601	1	-	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	601	NAG	O5-C1	2.41	1.47	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	NAG	C1-O5-C5	2.21	115.18	112.19

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	NAG	O5-C5-C6-O6
4	A	602	PEG	O1-C1-C2-O2
6	C	605	PG4	O2-C3-C4-O3
6	C	605	PG4	O3-C5-C6-O4
3	A	601	NAG	C4-C5-C6-O6
6	C	605	PG4	O1-C1-C2-O2
6	C	604	PG4	O2-C3-C4-O3
6	C	606	PG4	O1-C1-C2-O2
6	C	604	PG4	O4-C7-C8-O5
6	C	604	PG4	O3-C5-C6-O4
6	C	606	PG4	C4-C3-O2-C2
6	C	604	PG4	C5-C6-O4-C7
4	A	602	PEG	O2-C3-C4-O4
5	C	602	PGE	O1-C1-C2-O2
6	C	604	PG4	C6-C5-O3-C4
6	C	604	PG4	C8-C7-O4-C6
6	C	606	PG4	C1-C2-O2-C3
6	C	605	PG4	C3-C4-O3-C5
5	C	602	PGE	C1-C2-O2-C3
5	C	602	PGE	O3-C5-C6-O4
6	C	603	PG4	O2-C3-C4-O3
5	C	602	PGE	O2-C3-C4-O3
6	C	606	PG4	C3-C4-O3-C5
6	C	606	PG4	O2-C3-C4-O3

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Mol	Chain	Res	Type	Atoms
6	C	605	PG4	C5-C6-O4-C7
6	C	604	PG4	C3-C4-O3-C5

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	603	PG4	2	0
5	C	602	PGE	1	0
6	C	604	PG4	2	0
6	C	605	PG4	1	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	494/525 (94%)	0.08	22 (4%) 33 40	34, 46, 82, 113	5 (1%)
1	C	494/525 (94%)	0.15	24 (4%) 29 36	32, 47, 81, 116	1 (0%)
2	B	15/155 (9%)	0.42	2 (13%) 3 4	40, 49, 114, 123	0
2	D	15/155 (9%)	0.57	3 (20%) 1 1	39, 47, 105, 117	0
All	All	1018/1360 (74%)	0.12	51 (5%) 28 35	32, 47, 85, 123	6 (0%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	534	LEU	5.9
1	A	505	LEU	4.5
1	A	520	PRO	4.5
1	A	516	ARG	4.3
1	C	521	GLU	4.2
1	C	515	SER	4.2
1	C	288	ALA	4.2
1	A	515	SER	4.2
1	A	176	GLY	4.1
1	C	176	GLY	4.1
1	A	24	PRO	4.0
2	B	100	CYS	3.9
1	A	534	LEU	3.9
1	C	287	VAL	3.8
1	A	503	VAL	3.8
1	C	175	ARG	3.7
1	A	504	LEU	3.6
1	C	505	LEU	3.6
1	C	520	PRO	3.4
1	A	240	GLN	3.4
2	D	100	CYS	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	179	GLY	3.4
1	C	160	LEU	3.3
1	A	288	ALA	3.2
1	C	503	VAL	3.1
1	A	316	SER	3.1
2	D	102	SER	3.1
1	C	177	VAL	3.1
1	A	486	LEU	3.0
1	A	506	GLY	2.9
2	B	101	ASN	2.9
1	A	287	VAL	2.9
1	A	514	CYS	2.8
1	C	178	GLY	2.8
1	C	514	CYS	2.8
1	C	315	ALA	2.7
1	A	175	ARG	2.7
1	C	159	GLY	2.7
1	C	504	LEU	2.6
1	C	501	TRP	2.5
1	C	289	HIS	2.5
1	C	502	CYS	2.5
1	A	303	PRO	2.4
1	A	521	GLU	2.3
1	C	522	GLN	2.3
1	A	501	TRP	2.1
2	D	101	ASN	2.1
1	C	190	TRP	2.1
1	C	531	LEU	2.0
1	A	178	GLY	2.0
1	C	507	ARG	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](#)

There are no monosaccharides in this entry.

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. 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	B-factors(Å ²)	Q<0.9
6	PG4	C	606	13/13	0.65	0.33	53,78,89,91	0
6	PG4	C	605	13/13	0.81	0.21	55,69,77,80	0
6	PG4	C	604	13/13	0.82	0.18	63,72,84,84	0
4	PEG	A	602	7/7	0.83	0.21	60,63,71,79	0
5	PGE	C	602	10/10	0.84	0.18	52,61,72,74	0
6	PG4	C	603	13/13	0.91	0.16	45,57,64,67	0
3	NAG	A	601	14/15	0.94	0.12	46,53,70,75	0
3	NAG	C	601	14/15	0.96	0.11	41,50,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.