Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 07:20 pm BST

PDB ID	:	6EW9
Title	:	CRYSTAL STRUCTURE OF DEGS STRESS SENSOR PROTEASE IN COMPLEX WITH ACTIVATING DNRLGLVYQF PEPTIDE
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Deposited on	:	2017-11-03
Resolution	:	2.20 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 following versions of software and data (see references ⓘ) were used in the production of this report:

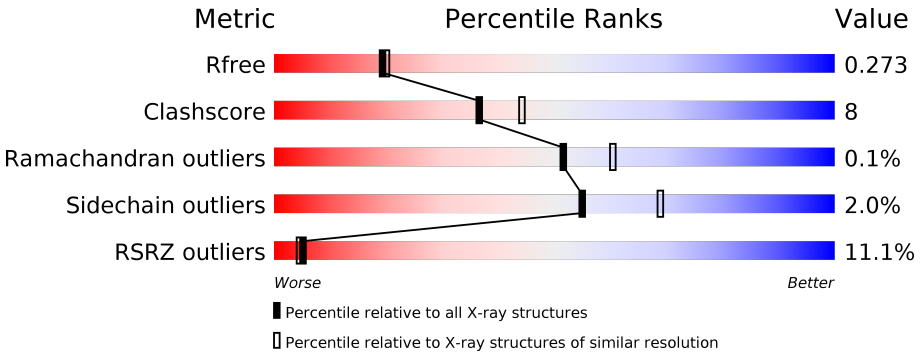
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 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 <=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	315	4% 84% 11%
1	B	315	15% 70% 23% 7%
1	C	315	10% 82% 13%
2	P	10	30% 20% 50%
2	Q	10	30% 30% 20% 50%
2	R	10	40% 10% 40% 50%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6919 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 Serine endoprotease DegS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	1	0
			2263	1417	402	439	5			
1	B	292	Total	C	N	O	S	0	0	0
			2169	1358	385	421	5			
1	C	302	Total	C	N	O	S	0	0	0
			2248	1404	403	436	5			

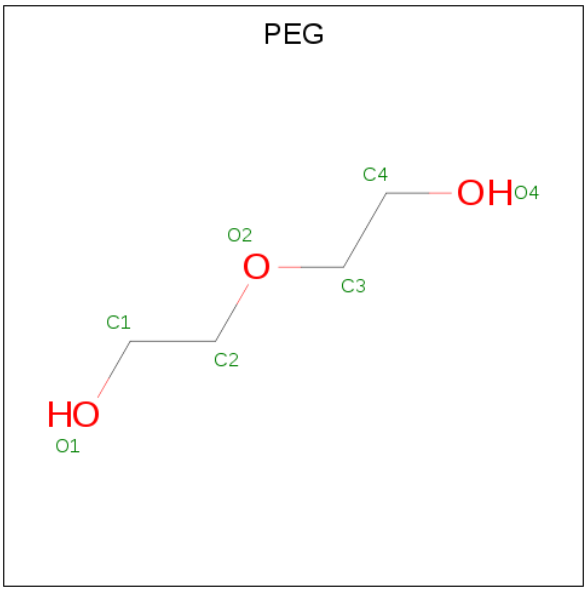
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	GLY	-	expression tag	UNP P0AEE3
A	42	PRO	-	expression tag	UNP P0AEE3
B	41	GLY	-	expression tag	UNP P0AEE3
B	42	PRO	-	expression tag	UNP P0AEE3
C	41	GLY	-	expression tag	UNP P0AEE3
C	42	PRO	-	expression tag	UNP P0AEE3

- Molecule 2 is a protein called DNRLGLVYQF PEPTIDE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	5	Total	C	N	O	0	0	0
			48	34	6	8			
2	Q	5	Total	C	N	O	0	0	0
			48	34	6	8			
2	R	5	Total	C	N	O	0	0	0
			48	34	6	8			

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



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

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	2	Total Ca 2 2	0	0
4	A	4	Total Ca 4 4	0	0
4	C	2	Total Ca 2 2	0	0

- Molecule 5 is water.

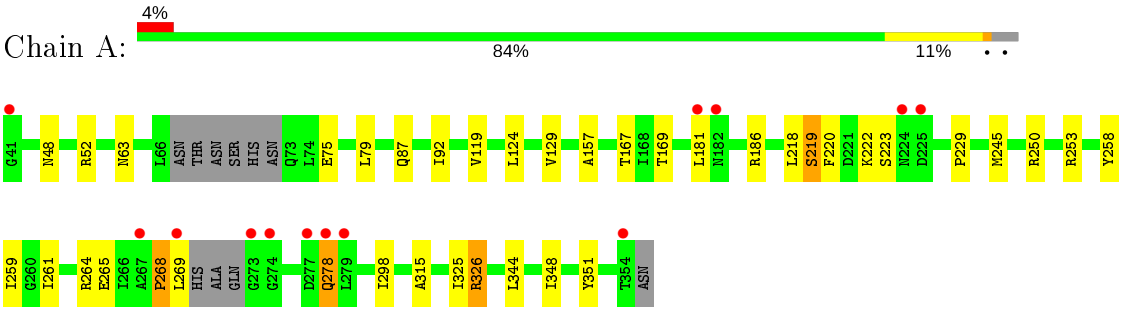
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	38	Total O 38 38	0	0
5	B	16	Total O 16 16	0	0
5	C	26	Total O 26 26	0	0

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 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.

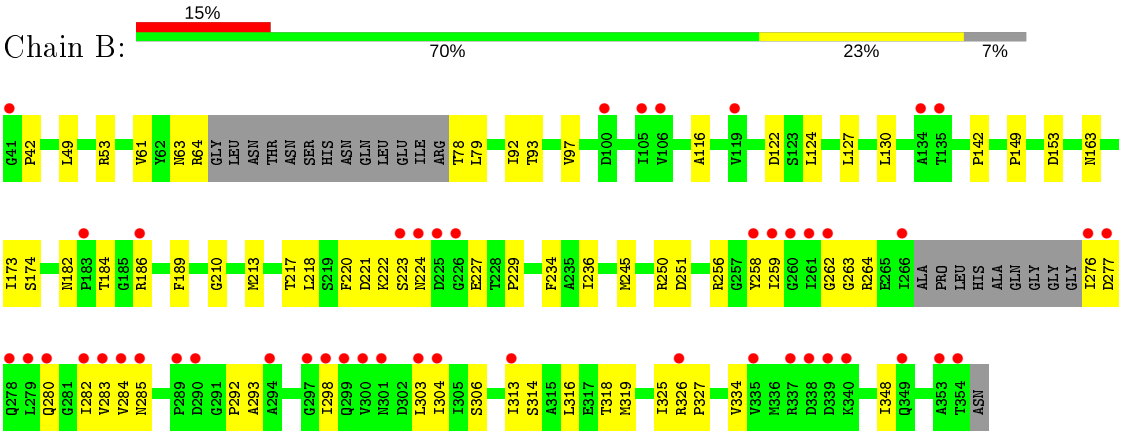
- Molecule 1: Serine endoprotease DegS

Chain A:



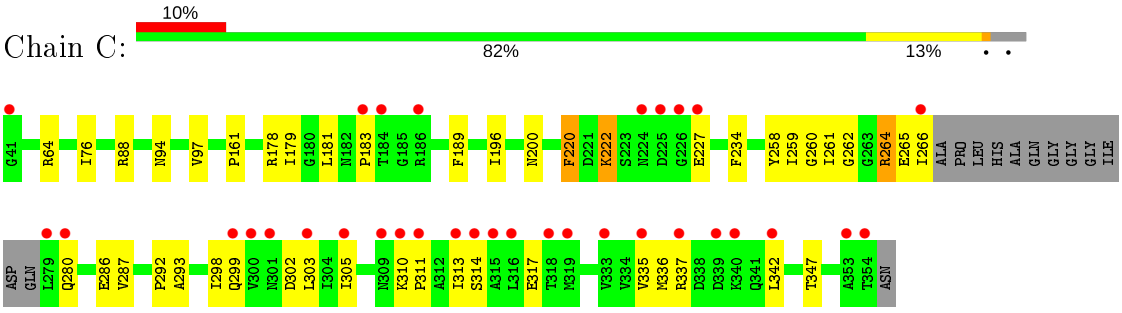
- Molecule 1: Serine endoprotease DegS

Chain B:



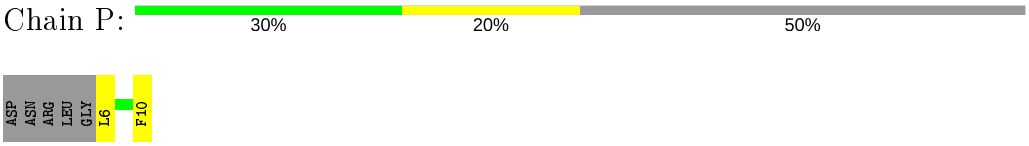
- Molecule 1: Serine endoprotease DegS

Chain C:



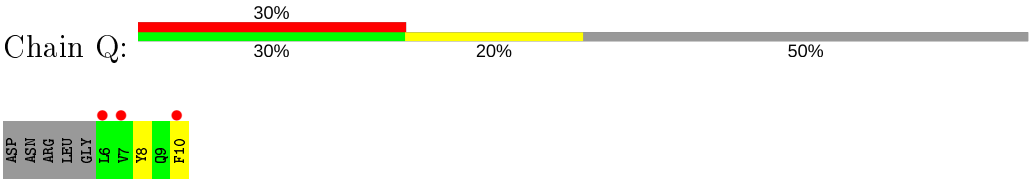
- Molecule 2: DNRLGLVYQF PEPTIDE

Chain P:



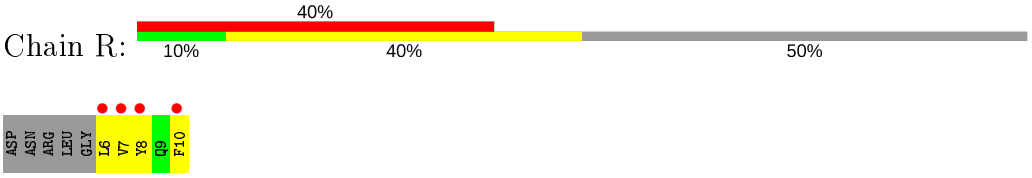
• Molecule 2: DNRLGLVYQF PEPTIDE

Chain Q:



• Molecule 2: DNRLGLVYQF PEPTIDE

Chain R:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	71.60Å 53.65Å 132.50Å 90.00° 101.80° 90.00°	Depositor
Resolution (Å)	43.38 – 2.20 43.38 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.6 (43.38-2.20) 99.7 (43.38-2.20)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.29 (at 2.20Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.226 , 0.273 0.226 , 0.273	Depositor DCC
R_{free} test set	2520 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	40.2	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 48.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6919	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% 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, 0.333 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: CA, PEG

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	0.27	0/2294	0.53	1/3120 (0.0%)
1	B	0.27	0/2196	0.52	0/2988
1	C	0.28	0/2277	0.53	2/3099 (0.1%)
2	P	0.28	0/49	0.51	0/64
2	Q	0.27	0/49	0.42	0/64
2	R	0.26	0/49	0.35	0/64
All	All	0.27	0/6914	0.53	3/9399 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	227	GLU	C-N-CA	-7.46	103.06	121.70
1	C	181	LEU	C-N-CA	-5.38	108.25	121.70
1	A	326	ARG	NE-CZ-NH1	5.35	122.98	120.30

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	2263	0	2313	25	0
1	B	2169	0	2214	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2248	0	2292	36	0
2	P	48	0	45	1	0
2	Q	48	0	45	3	0
2	R	48	0	45	5	0
3	A	7	0	10	1	0
4	A	4	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
5	A	38	0	0	2	0
5	B	16	0	0	0	0
5	C	26	0	0	0	0
All	All	6919	0	6964	109	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 8.

All (109) 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:92:ILE:HD11	1:A:245:MET:HB2	1.53	0.88
1:C:220:PHE:HE2	1:C:222:LYS:HG2	1.46	0.80
1:B:284:VAL:HG23	1:B:304:ILE:HD11	1.65	0.79
1:A:181:LEU:HD13	1:A:222:LYS:HD2	1.69	0.73
1:B:293:ALA:HB1	1:B:298:ILE:HD11	1.73	0.70
1:C:313:ILE:HD12	1:C:314:SER:N	2.07	0.69
1:B:127:LEU:HD11	1:B:236:ILE:HD13	1.76	0.67
1:B:280:GLN:HB2	1:B:313:ILE:HG12	1.77	0.66
1:C:261:ILE:HG22	1:C:287:VAL:HG22	1.76	0.66
1:A:219:SER:OG	5:A:501:HOH:O	2.15	0.65
1:B:124:LEU:HD22	1:B:186:ARG:HH12	1.60	0.64
1:B:250:ARG:NH1	1:B:251:ASP:OD1	2.32	0.62
1:C:303:LEU:HD12	1:C:305:ILE:HG23	1.81	0.62
1:C:259:ILE:HG22	1:C:293:ALA:HB2	1.82	0.62
1:C:264:ARG:HD3	1:C:265:GLU:H	1.65	0.61
1:A:220:PHE:HD2	1:A:229:PRO:HG3	1.66	0.61
1:A:181:LEU:HD21	1:A:220:PHE:CD1	2.37	0.60
1:C:303:LEU:CD1	1:C:305:ILE:HG23	2.32	0.60
1:B:319:MET:HG3	2:Q:10:PHE:HB3	1.83	0.59
1:B:220:PHE:CZ	1:B:223:SER:HB3	2.38	0.59
1:A:269:LEU:HD13	3:A:401:PEG:H22	1.85	0.58
1:B:122:ASP:OD1	1:B:256:ARG:NH2	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:310:LYS:HG2	1:C:311:PRO:HD2	1.85	0.57
1:B:186:ARG:HE	1:B:319:MET:HG2	1.70	0.56
1:C:264:ARG:HD3	1:C:265:GLU:N	2.21	0.55
1:C:258:TYR:CZ	1:C:260:GLY:HA2	2.42	0.54
1:C:261:ILE:HD11	2:R:10:PHE:CZ	2.42	0.54
1:A:63:ASN:HB2	1:A:79:LEU:HD11	1.91	0.52
1:C:293:ALA:HB1	1:C:298:ILE:HG21	1.91	0.52
1:A:92:ILE:HG12	1:A:245:MET:HE3	1.91	0.52
1:A:265:GLU:HG3	1:A:315:ALA:HB2	1.92	0.51
1:B:122:ASP:CG	1:B:256:ARG:HH22	2.13	0.51
1:A:326:ARG:HG3	1:A:326:ARG:HH11	1.75	0.51
1:B:61:VAL:HG23	1:B:79:LEU:HB2	1.92	0.51
1:C:313:ILE:HD12	1:C:314:SER:H	1.72	0.51
1:C:293:ALA:HB1	1:C:298:ILE:CG2	2.41	0.51
1:C:335:VAL:HG12	1:C:342:LEU:HB2	1.92	0.51
1:B:49:LEU:HD21	1:B:53:ARG:CZ	2.42	0.50
1:A:259:ILE:HG23	1:A:261:ILE:HG23	1.94	0.50
1:C:94:ASN:HB2	1:C:97:VAL:HG23	1.94	0.49
1:B:92:ILE:HD11	1:B:245:MET:HB2	1.94	0.49
1:C:189:PHE:CG	1:C:234:PHE:HB3	2.48	0.49
1:A:167:THR:HG23	1:B:174:SER:HB3	1.94	0.49
1:A:298:ILE:HG13	1:A:344:LEU:HD12	1.95	0.49
1:C:262:GLY:HA2	2:R:8:TYR:O	2.13	0.49
1:A:268:PRO:HB3	1:A:278:GLN:HB3	1.95	0.48
1:C:302:ASP:HB3	1:C:335:VAL:CG2	2.43	0.48
1:B:264:ARG:HE	1:B:285:ASN:ND2	2.11	0.48
1:C:311:PRO:HB2	1:C:313:ILE:HG23	1.96	0.48
1:A:258:TYR:H	1:A:348:ILE:HG23	1.79	0.48
1:B:326:ARG:HG3	1:B:327:PRO:HD2	1.95	0.47
1:A:261:ILE:HD11	2:P:10:PHE:CZ	2.48	0.47
1:C:264:ARG:HG2	2:R:6:LEU:O	2.13	0.47
1:B:153:ASP:HB2	1:B:173:ILE:HD12	1.96	0.47
1:B:262:GLY:HA2	2:Q:8:TYR:O	2.15	0.47
1:C:88:ARG:HH11	1:C:88:ARG:HG3	1.80	0.47
1:B:264:ARG:O	1:B:282:ILE:HA	2.14	0.46
1:C:292:PRO:HB3	1:C:347:THR:O	2.15	0.46
1:B:263:GLY:HA2	1:B:283:VAL:O	2.14	0.46
1:B:306:SER:HB2	1:B:334:VAL:HB	1.96	0.46
1:C:64:ARG:HG2	1:C:76:ILE:HD13	1.97	0.46
1:B:93:THR:OG1	1:B:97:VAL:HG11	2.15	0.46
1:C:179:ILE:HD12	1:C:183:PRO:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:335:VAL:HG22	1:C:336:MET:H	1.81	0.46
1:B:218:LEU:HD23	1:B:234:PHE:HB2	1.97	0.46
1:C:313:ILE:HD11	1:C:317:GLU:HB2	1.97	0.45
1:B:325:ILE:HG21	1:B:348:ILE:HG13	1.99	0.45
1:B:319:MET:HE1	2:Q:8:TYR:CD2	2.52	0.45
1:A:48:ASN:O	1:A:52:ARG:HG3	2.17	0.45
1:B:217:THR:OG1	1:B:236:ILE:HD12	2.16	0.45
1:C:261:ILE:HD11	2:R:10:PHE:CE2	2.51	0.45
1:A:157:ALA:HB3	1:A:169:THR:OG1	2.16	0.45
1:B:224:ASN:O	1:B:227:GLU:HG2	2.17	0.45
1:B:124:LEU:HD22	1:B:186:ARG:NH1	2.30	0.45
1:C:293:ALA:O	1:C:298:ILE:HG22	2.17	0.44
1:B:221:ASP:O	1:B:222:LYS:HD2	2.17	0.44
1:B:293:ALA:O	1:B:298:ILE:HG12	2.16	0.44
2:R:6:LEU:HB3	2:R:7:VAL:H	1.68	0.44
1:C:178:ARG:O	1:C:189:PHE:N	2.47	0.44
1:A:186:ARG:NH1	5:A:503:HOH:O	2.29	0.44
1:A:325:ILE:HG21	1:A:348:ILE:HG13	2.00	0.44
1:B:189:PHE:CG	1:B:234:PHE:HB3	2.53	0.44
1:B:259:ILE:O	1:B:292:PRO:HG2	2.18	0.43
1:B:78:THR:HG21	1:B:163:ASN:HB2	1.99	0.43
1:B:149:PRO:HB3	1:B:213:MET:HE3	2.01	0.43
1:B:223:SER:HB2	1:B:224:ASN:H	1.55	0.43
1:A:181:LEU:HD12	1:A:218:LEU:HD21	2.00	0.43
1:B:182:ASN:ND2	1:B:184:THR:O	2.52	0.43
1:B:227:GLU:O	1:B:229:PRO:HD3	2.19	0.43
1:C:266:ILE:HB	1:C:280:GLN:HG2	2.00	0.43
1:B:142:PRO:HG2	1:B:210:GLY:O	2.18	0.43
1:B:63:ASN:HB2	1:B:79:LEU:HD11	2.01	0.42
1:C:302:ASP:OD2	1:C:337:ARG:HB3	2.19	0.42
1:B:264:ARG:HE	1:B:285:ASN:HD21	1.67	0.42
1:B:280:GLN:H	1:B:280:GLN:HG2	1.64	0.42
1:B:127:LEU:CD1	1:B:236:ILE:HD13	2.45	0.42
1:C:299:GLN:HA	1:C:299:GLN:NE2	2.34	0.42
1:B:276:ILE:O	1:B:277:ASP:HB2	2.20	0.42
1:B:314:SER:O	1:B:318:THR:HG23	2.20	0.42
1:B:64:ARG:HH11	1:B:64:ARG:HG3	1.86	0.41
1:A:181:LEU:HD23	1:A:181:LEU:HA	1.75	0.41
1:C:196:ILE:HD13	1:C:196:ILE:HA	1.87	0.41
1:A:181:LEU:HD11	1:A:220:PHE:HA	2.03	0.41
1:C:161:PRO:HG3	1:C:200:ASN:OD1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:VAL:HB	1:A:129:VAL:HG12	2.02	0.41
1:C:286:GLU:HG2	1:C:287:VAL:N	2.35	0.41
1:A:124:LEU:HD11	1:A:351:TYR:CE1	2.56	0.40
1:B:116:ALA:HB1	1:B:130:LEU:HB3	2.03	0.40
1:B:282:ILE:O	1:B:303:LEU:HD12	2.22	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	300/315 (95%)	290 (97%)	9 (3%)	1 (0%)	41	46
1	B	286/315 (91%)	276 (96%)	10 (4%)	0	100	100
1	C	298/315 (95%)	289 (97%)	9 (3%)	0	100	100
2	P	3/10 (30%)	3 (100%)	0	0	100	100
2	Q	3/10 (30%)	3 (100%)	0	0	100	100
2	R	3/10 (30%)	3 (100%)	0	0	100	100
All	All	893/975 (92%)	864 (97%)	28 (3%)	1 (0%)	51	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	PRO

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	245/253 (97%)	237 (97%)	8 (3%)	38 49
1	B	236/253 (93%)	233 (99%)	3 (1%)	69 81
1	C	245/253 (97%)	242 (99%)	3 (1%)	71 83
2	P	5/9 (56%)	4 (80%)	1 (20%)	1 1
2	Q	5/9 (56%)	5 (100%)	0	100 100
2	R	5/9 (56%)	5 (100%)	0	100 100
All	All	741/786 (94%)	726 (98%)	15 (2%)	55 69

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	GLU
1	A	87	GLN
1	A	219	SER
1	A	223	SER
1	A	250	ARG
1	A	253	ARG
1	A	264	ARG
1	A	278	GLN
1	B	42	PRO
1	B	258	TYR
1	B	316	LEU
1	C	220	PHE
1	C	222	LYS
1	C	264	ARG
2	P	6	LEU

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

Mol	Chain	Res	Type
1	C	299	GLN

5.3.3 RNA

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 9 ligands modelled in this entry, 8 are monoatomic - leaving 1 for Mogul analysis.

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	PEG	A	401	-	6,6,6	0.38	0	5,5,5	0.38	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
3	PEG	A	401	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	PEG	C4-C3-O2-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	PEG	1	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	305/315 (96%)	0.39	13 (4%) 35 33	29, 42, 73, 91	0
1	B	292/315 (92%)	1.03	48 (16%) 1 1	36, 57, 83, 94	0
1	C	302/315 (95%)	0.67	33 (10%) 5 5	28, 43, 85, 92	0
2	P	5/10 (50%)	1.17	0 100 100	51, 56, 67, 76	0
2	Q	5/10 (50%)	2.83	3 (60%) 0 0	70, 72, 77, 77	0
2	R	5/10 (50%)	2.94	4 (80%) 0 0	76, 77, 83, 83	0
All	All	914/975 (93%)	0.72	101 (11%) 5 4	28, 46, 83, 94	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	183	PRO	9.9
1	B	226	GLY	8.1
1	B	279	LEU	8.1
1	C	279	LEU	8.1
1	C	266	ILE	6.5
1	B	224	ASN	6.0
1	C	225	ASP	5.6
1	C	313	ILE	5.5
2	Q	6	LEU	5.3
1	C	184	THR	5.3
1	C	280	GLN	5.2
1	B	313	ILE	4.9
1	B	340	LYS	4.8
1	B	266	ILE	4.8
1	C	226	GLY	4.7
1	A	224	ASN	4.7
2	R	7	VAL	4.6
1	C	224	ASN	4.6
1	B	304	ILE	4.4

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Mol	Chain	Res	Type	RSRZ
1	B	261	ILE	4.3
1	B	289	PRO	4.3
1	B	294	ALA	4.3
1	B	276	ILE	4.2
2	Q	10	PHE	4.2
1	B	303	LEU	4.2
1	B	280	GLN	4.1
1	C	339	ASP	4.1
1	A	41	GLY	4.1
1	B	223	SER	4.0
2	R	8	TYR	3.9
1	B	300	VAL	3.8
2	R	10	PHE	3.8
1	B	225	ASP	3.8
1	C	303	LEU	3.7
1	C	305	ILE	3.7
1	B	301	ASN	3.7
1	A	273	GLY	3.6
1	B	183	PRO	3.5
1	A	274	GLY	3.5
1	A	279	LEU	3.3
1	B	135	THR	3.3
1	A	269	LEU	3.3
1	B	282	ILE	3.2
1	C	353	ALA	3.2
1	B	258	TYR	3.2
1	C	340	LYS	3.1
1	B	186	ARG	3.1
1	B	134	ALA	3.1
1	C	342	LEU	3.1
1	B	335	VAL	3.1
1	B	353	ALA	3.1
1	A	181	LEU	3.0
1	A	225	ASP	3.0
1	B	106	VAL	3.0
1	C	335	VAL	3.0
1	A	182	ASN	3.0
1	B	285	ASN	2.9
1	C	300	VAL	2.9
2	Q	7	VAL	2.8
1	B	354	THR	2.8
1	C	309	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	41	GLY	2.7
1	B	259	ILE	2.7
1	C	186	ARG	2.7
1	C	337	ARG	2.7
1	B	277	ASP	2.7
2	R	6	LEU	2.7
1	B	262	GLY	2.7
1	C	311	PRO	2.6
1	A	278	GLN	2.6
1	B	284	VAL	2.5
1	A	277	ASP	2.5
1	C	299	GLN	2.5
1	A	354	THR	2.5
1	B	298	ILE	2.5
1	C	333	VAL	2.4
1	C	301	ASN	2.4
1	B	119	VAL	2.4
1	C	354	THR	2.4
1	B	326	ARG	2.4
1	B	283	VAL	2.4
1	B	339	ASP	2.3
1	C	319	MET	2.3
1	B	290	ASP	2.3
1	C	318	THR	2.2
1	B	105	ILE	2.2
1	B	349	GLN	2.2
1	B	100	ASP	2.2
1	B	260	GLY	2.2
1	B	297	GLY	2.2
1	B	299	GLN	2.1
1	C	314	SER	2.1
1	C	316	LEU	2.1
1	B	278	GLN	2.1
1	C	227	GLU	2.1
1	A	267	ALA	2.1
1	C	310	LYS	2.0
1	C	41	GLY	2.0
1	B	337	ARG	2.0
1	B	338	ASP	2.0
1	C	315	ALA	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 carbohydrates 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, 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	B-factors(Å^2)	Q<0.9
4	CA	C	402	1/1	0.49	0.34	68,68,68,68	0
4	CA	B	401	1/1	0.52	0.16	70,70,70,70	0
3	PEG	A	401	7/7	0.78	0.31	80,82,85,88	3
4	CA	C	401	1/1	0.83	0.13	72,72,72,72	0
4	CA	A	404	1/1	0.86	0.08	80,80,80,80	0
4	CA	A	403	1/1	0.88	0.12	73,73,73,73	0
4	CA	B	402	1/1	0.89	0.08	70,70,70,70	0
4	CA	A	405	1/1	0.94	0.17	80,80,80,80	0
4	CA	A	402	1/1	0.98	0.16	77,77,77,77	0

6.5 Other polymers (i)

There are no such residues in this entry.