Full wwPDB X-ray Structure Validation Report (i)

Dec 17, 2023 – 06:26 PM EST

PDB ID	:	4WMO
Title	:	Selenomethionine derivative of Xenopus laevis embryonic epidermal lectin carbohydrate-binding domain
Authors	:	Wangkanont, K.; Kiessling, L.L.; Forest, K.T.
Deposited on	:	2014-10-09
Resolution	:	2.30 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 (i) 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 (i)) were used in the production of this report:

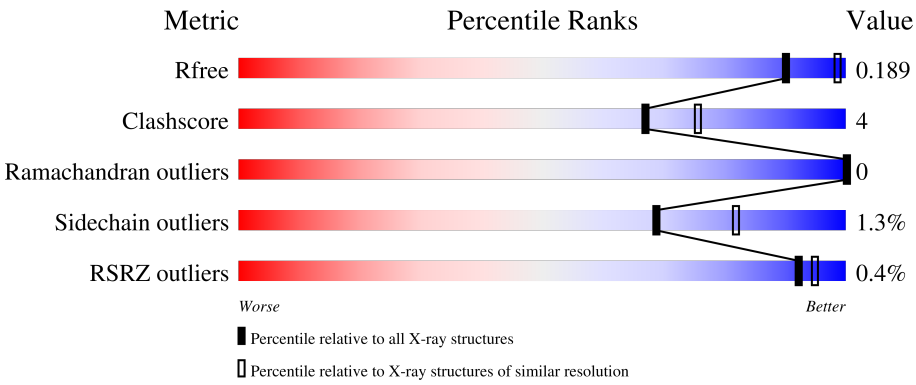
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA,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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.30 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 <=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	289	87% 9% •
1	B	289	88% 7% • •
1	C	289	87% 8% •
1	D	289	.% 89% 7% •
1	E	289	.% 89% 7% •

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Mol	Chain	Length	Quality of chain
1	F	289	87% 9% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 14639 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.

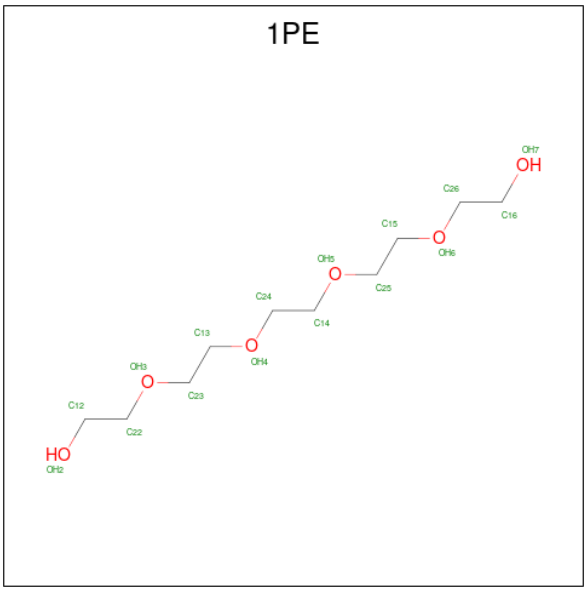
- Molecule 1 is a protein called XEEL protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	F	277	2153	1350	367	421	8	7	0	2	0
1	C	277	2153	1350	367	421	8	7	0	2	0
1	B	277	2158	1353	370	420	8	7	0	2	0
1	D	277	2147	1347	366	419	8	7	0	1	0
1	E	277	2147	1347	366	419	8	7	0	1	0
1	A	277	2169	1359	374	421	8	7	0	3	0

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
2	F	3	3	3	0	0
2	C	3	3	3	0	0
2	B	3	3	3	0	0
2	D	3	3	3	0	0
2	E	3	3	3	0	0
2	A	3	3	3	0	0

- Molecule 3 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	F	1	16	10	6	0	0
			Total	C	O		
3	F	1	16	10	6	0	0
			Total	C	O		
3	C	1	16	10	6	0	0
			Total	C	O		
3	C	1	16	10	6	0	0
			Total	C	O		
3	C	1	16	10	6	0	0
			Total	C	O		
3	B	1	16	10	6	0	0
			Total	C	O		
3	B	1	16	10	6	0	0
			Total	C	O		
3	D	1	16	10	6	0	0
			Total	C	O		
3	D	1	16	10	6	0	0
			Total	C	O		
3	E	1	16	10	6	0	0
			Total	C	O		
3	E	1	16	10	6	0	0
			Total	C	O		
3	A	1	16	10	6	0	0
			Total	C	O		
3	A	1	16	10	6	0	0
			Total	C	O		
3	A	1	16	10	6	0	0

- Molecule 4 is water.

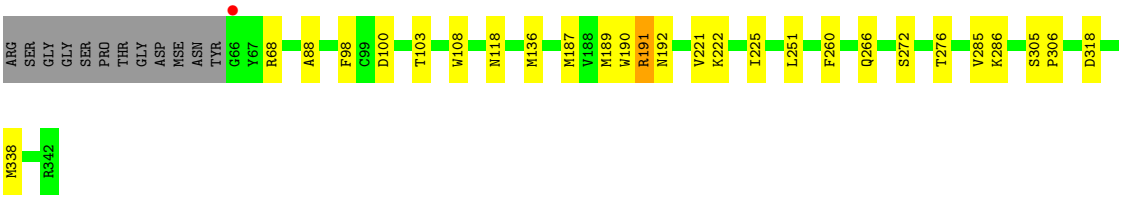
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	F	236	236	236	0	0
			Total	O		
4	C	228	228	228	0	0
			Total	O		
4	B	236	236	236	0	0
			Total	O		
4	D	262	262	262	0	0
			Total	O		
4	E	241	241	241	0	0
			Total	O		
4	A	267	267	267	0	0

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.

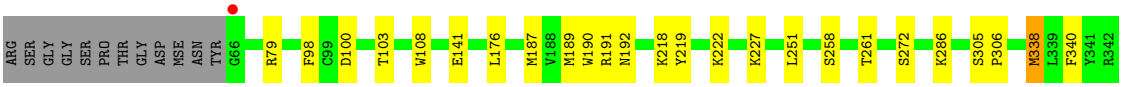
- Molecule 1: XEEL protein

Chain F:



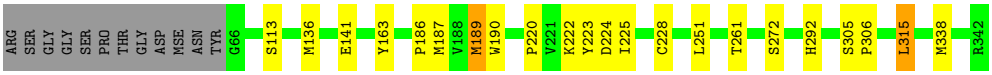
- Molecule 1: XEEL protein

Chain C:



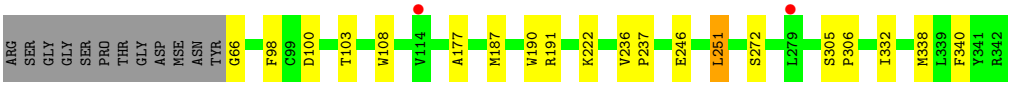
- Molecule 1: XEEL protein

Chain B:



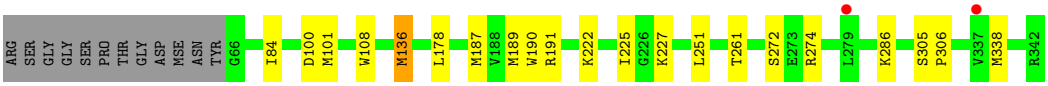
- Molecule 1: XEEL protein

Chain D:



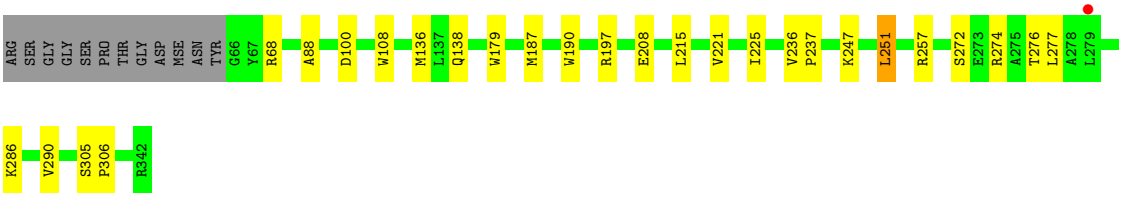
- Molecule 1: XEEL protein

Chain E:



• Molecule 1: XEEL protein

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	123.65Å 111.07Å 123.61Å 90.00° 119.72° 90.00°	Depositor
Resolution (Å)	30.09 – 2.30 30.09 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.09-2.30) 99.8 (30.09-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.17	Depositor
$< I/\sigma(I) >$ [1]	2.75 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.160 , 0.188 0.161 , 0.189	Depositor DCC
R_{free} test set	6448 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	19.4	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 56.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h-l 0.000 for -h-l,k,h 0.000 for -h-l,-k,l 0.015 for h,-k,-h-l 0.000 for l,-k,h	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	14639	wwPDB-VP
Average B, all atoms (Å^2)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.74 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.4617e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 1PE

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	0.46	0/2219	0.61	2/2998 (0.1%)
1	B	0.46	0/2208	0.63	4/2984 (0.1%)
1	C	0.45	0/2203	0.61	2/2978 (0.1%)
1	D	0.47	0/2197	0.62	1/2970 (0.0%)
1	E	0.45	0/2197	0.62	4/2970 (0.1%)
1	F	0.44	0/2203	0.60	3/2978 (0.1%)
All	All	0.45	0/13227	0.61	16/17878 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	136[A]	MSE	CG-SE-CE	6.37	112.90	98.90
1	B	136[B]	MSE	CG-SE-CE	6.37	112.90	98.90
1	B	189	MSE	CG-SE-CE	6.23	112.61	98.90
1	F	189	MSE	CG-SE-CE	6.05	112.22	98.90
1	E	101	MSE	CG-SE-CE	6.03	112.17	98.90
1	F	136[A]	MSE	CG-SE-CE	5.66	111.35	98.90
1	F	136[B]	MSE	CG-SE-CE	5.66	111.35	98.90
1	A	136[A]	MSE	CG-SE-CE	5.60	111.22	98.90
1	A	136[B]	MSE	CG-SE-CE	5.60	111.22	98.90
1	E	136[A]	MSE	CG-SE-CE	5.48	110.96	98.90
1	E	136[B]	MSE	CG-SE-CE	5.48	110.96	98.90
1	B	338	MSE	CA-CB-CG	-5.47	104.00	113.30
1	E	189	MSE	CG-SE-CE	5.46	110.91	98.90
1	C	189	MSE	CG-SE-CE	5.17	110.27	98.90
1	D	338	MSE	CG-SE-CE	5.09	110.09	98.90
1	C	338	MSE	CA-CB-CG	-5.03	104.75	113.30

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	2169	0	2031	18	0
1	B	2158	0	2019	14	0
1	C	2153	0	2011	17	0
1	D	2147	0	2007	12	0
1	E	2147	0	2007	20	0
1	F	2153	0	2011	19	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
2	E	3	0	0	0	0
2	F	3	0	0	0	0
3	A	48	0	66	5	0
3	B	32	0	44	7	0
3	C	48	0	66	5	0
3	D	32	0	44	1	0
3	E	32	0	44	10	0
3	F	32	0	44	7	0
4	A	267	0	0	3	0
4	B	236	0	0	0	0
4	C	228	0	0	3	0
4	D	262	0	0	1	0
4	E	241	0	0	3	0
4	F	236	0	0	2	0
All	All	14639	0	12394	99	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 4.

All (99) 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:E:286:LYS:HZ3	3:E:405:1PE:H261	1.52	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:CYS:HB3	3:B:405:1PE:H162	1.73	0.71
1:A:225:ILE:HD12	3:A:404:1PE:H131	1.74	0.69
1:D:100:ASP:HB3	1:D:108:TRP:HB2	1.76	0.68
1:B:222:LYS:HZ2	3:B:404:1PE:H151	1.59	0.67
1:D:222:LYS:HE2	3:D:404:1PE:H132	1.76	0.67
1:F:192:ASN:OD1	4:F:627:HOH:O	2.12	0.67
1:C:222:LYS:HZ3	3:C:404:1PE:H142	1.60	0.66
1:E:261:THR:HB	3:E:405:1PE:H242	1.77	0.66
1:D:66:GLY:N	4:D:687:HOH:O	2.29	0.65
1:F:222:LYS:HZ2	3:F:404:1PE:H242	1.62	0.65
1:F:286:LYS:NZ	3:F:405:1PE:H231	2.10	0.65
1:F:192:ASN:ND2	4:F:725:HOH:O	2.25	0.64
1:A:100:ASP:HB3	1:A:108:TRP:HB2	1.80	0.64
1:C:103:THR:OG1	1:C:191:ARG:NH1	2.33	0.61
1:A:286:LYS:HD3	3:A:405:1PE:H251	1.83	0.60
1:C:286:LYS:HE2	3:C:405:1PE:H122	1.83	0.59
1:C:100:ASP:HB3	1:C:108:TRP:HB2	1.85	0.59
1:F:100:ASP:HB3	1:F:108:TRP:HB2	1.84	0.59
1:E:225:ILE:HD11	3:E:404:1PE:H251	1.84	0.58
1:E:286:LYS:NZ	3:E:405:1PE:H261	2.20	0.57
1:F:286:LYS:NZ	3:F:405:1PE:H152	2.20	0.57
1:B:141:GLU:OE2	3:B:404:1PE:H252	2.05	0.57
1:B:222:LYS:NZ	3:B:404:1PE:H151	2.20	0.56
1:A:187:MSE:HA	1:A:190:TRP:CE2	2.41	0.55
1:F:225:ILE:CD1	3:F:404:1PE:H151	2.38	0.54
1:F:286:LYS:HE2	3:F:405:1PE:H222	1.88	0.54
1:D:187:MSE:HA	1:D:190:TRP:CE2	2.44	0.53
1:E:190:TRP:CH2	1:E:338:MSE:HE1	2.43	0.53
1:F:103:THR:OG1	1:F:191:ARG:NH1	2.41	0.53
1:F:286:LYS:HZ2	3:F:405:1PE:H231	1.71	0.53
1:F:318:ASP:OD1	1:C:79:ARG:NH1	2.41	0.52
1:E:191:ARG:NH1	4:E:683:HOH:O	2.29	0.52
1:E:286:LYS:HZ3	3:E:405:1PE:H251	1.74	0.52
1:F:222:LYS:NZ	3:F:404:1PE:H242	2.24	0.52
1:C:227:LYS:HD2	3:C:406:1PE:H221	1.91	0.52
1:C:261:THR:HB	3:C:405:1PE:H242	1.92	0.52
1:A:286:LYS:HZ3	3:A:405:1PE:H261	1.76	0.51
1:B:222:LYS:HZ3	3:B:404:1PE:H121	1.76	0.50
1:E:222:LYS:HZ3	3:E:404:1PE:H152	1.75	0.50
1:E:225:ILE:CD1	3:E:404:1PE:H251	2.41	0.50
1:D:177:ALA:HB3	1:D:340:PHE:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:286:LYS:HD3	3:E:405:1PE:H251	1.94	0.49
1:D:246:GLU:H	1:D:246:GLU:CD	2.15	0.48
1:B:292:HIS:O	1:B:315:LEU:HD22	2.13	0.48
1:C:218:LYS:HD2	1:C:219:TYR:CE2	2.48	0.48
1:F:118:ASN:HD21	1:E:136[B]:MSE:SE	2.47	0.48
1:A:274:ARG:NH1	4:A:502:HOH:O	2.46	0.47
1:E:225:ILE:HG22	4:E:586:HOH:O	2.14	0.47
1:C:192:ASN:ND2	4:C:666:HOH:O	2.48	0.47
1:C:187:MSE:HA	1:C:190:TRP:CE2	2.50	0.46
1:A:68:ARG:HG2	1:A:88:ALA:O	2.15	0.46
1:C:305:SER:HA	1:C:306:PRO:HA	1.75	0.46
1:F:98:PHE:HB2	1:F:187:MSE:HE2	1.97	0.46
1:F:266:GLN:OE1	1:F:285:VAL:HG22	2.16	0.46
1:C:98:PHE:HB2	1:C:187:MSE:HE2	1.99	0.45
1:B:220:PRO:HB2	1:B:225:ILE:HD12	1.98	0.45
1:B:113:SER:HB3	1:B:163:TYR:HB3	1.99	0.45
1:D:103:THR:OG1	1:D:191:ARG:NH2	2.49	0.44
1:A:277:LEU:HD23	1:A:290:VAL:HG23	1.99	0.44
1:A:286:LYS:NZ	3:A:405:1PE:H261	2.31	0.44
1:F:68:ARG:HG2	1:F:88:ALA:O	2.18	0.44
1:E:100:ASP:HB3	1:E:108:TRP:HB2	1.99	0.44
1:F:305:SER:HA	1:F:306:PRO:HA	1.76	0.44
1:E:305:SER:HA	1:E:306:PRO:HA	1.85	0.44
1:A:179:TRP:CG	1:A:197:ARG:HG2	2.53	0.44
1:C:191:ARG:NE	4:C:689:HOH:O	2.38	0.44
1:B:187:MSE:HA	1:B:190:TRP:CE2	2.52	0.44
1:C:190:TRP:CZ3	1:C:338:MSE:HE1	2.53	0.44
1:E:227:LYS:HE2	1:E:274:ARG:NH2	2.33	0.44
1:E:191:ARG:HD3	4:E:683:HOH:O	2.19	0.43
1:F:221:VAL:HA	1:F:276:THR:HG21	2.00	0.43
1:E:187:MSE:HA	1:E:190:TRP:CE2	2.54	0.43
1:A:208:GLU:HG3	1:A:215:LEU:HD12	2.00	0.43
1:C:141:GLU:HB2	4:C:554:HOH:O	2.19	0.42
1:F:190:TRP:CH2	1:F:338:MSE:HE1	2.54	0.42
1:E:84:ILE:HD11	1:E:187:MSE:HE1	2.01	0.42
1:A:221:VAL:HA	1:A:276:THR:HG21	2.01	0.42
1:A:225:ILE:HG22	4:A:548:HOH:O	2.18	0.42
1:A:247:LYS:HG2	1:A:251:LEU:HD22	2.01	0.42
1:A:286:LYS:CE	3:A:405:1PE:H261	2.49	0.42
1:B:305:SER:HA	1:B:306:PRO:HA	1.86	0.42
1:C:176:LEU:HD12	1:C:340:PHE:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:PRO:HD2	1:B:189:MSE:HE2	2.02	0.42
1:B:223:TYR:CE2	1:B:224:ASP:HB2	2.54	0.42
1:D:98:PHE:HB2	1:D:187:MSE:HE2	2.01	0.42
1:E:178:LEU:HA	1:E:338:MSE:O	2.20	0.42
1:E:222:LYS:HZ3	3:E:404:1PE:C15	2.34	0.41
1:B:222:LYS:NZ	3:B:404:1PE:H121	2.35	0.41
1:A:257[B]:ARG:NH2	4:A:691:HOH:O	2.22	0.41
1:C:258:SER:O	3:C:405:1PE:H231	2.20	0.41
3:E:405:1PE:H251	3:E:405:1PE:H241	1.88	0.41
1:D:251:LEU:HB3	1:D:332:ILE:HD12	2.01	0.41
1:B:222:LYS:HE3	3:B:404:1PE:H262	2.03	0.41
1:A:305:SER:HA	1:A:306:PRO:HA	1.83	0.41
1:D:187:MSE:HA	1:D:190:TRP:CD2	2.56	0.40
1:A:236:VAL:HA	1:A:237:PRO:HD3	1.99	0.40
1:D:305:SER:HA	1:D:306:PRO:HA	1.85	0.40
1:D:236:VAL:HA	1:D:237:PRO:HD3	1.99	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	278/289 (96%)	269 (97%)	9 (3%)	0	100	100
1	B	277/289 (96%)	268 (97%)	9 (3%)	0	100	100
1	C	277/289 (96%)	265 (96%)	12 (4%)	0	100	100
1	D	276/289 (96%)	263 (95%)	13 (5%)	0	100	100
1	E	276/289 (96%)	266 (96%)	10 (4%)	0	100	100
1	F	277/289 (96%)	266 (96%)	11 (4%)	0	100	100
All	All	1661/1734 (96%)	1597 (96%)	64 (4%)	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	226/225 (100%)	223 (99%)	3 (1%)	69	82
1	B	225/225 (100%)	221 (98%)	4 (2%)	59	75
1	C	225/225 (100%)	223 (99%)	2 (1%)	78	89
1	D	224/225 (100%)	222 (99%)	2 (1%)	78	89
1	E	224/225 (100%)	222 (99%)	2 (1%)	78	89
1	F	225/225 (100%)	221 (98%)	4 (2%)	59	75
All	All	1349/1350 (100%)	1332 (99%)	17 (1%)	69	82

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	191	ARG
1	F	251	LEU
1	F	260	PHE
1	F	272	SER
1	C	251	LEU
1	C	272	SER
1	B	251	LEU
1	B	261	THR
1	B	272	SER
1	B	315	LEU
1	D	251	LEU
1	D	272	SER
1	E	251	LEU
1	E	272	SER
1	A	138	GLN
1	A	251	LEU
1	A	272	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are

no such sidechains identified.

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)

Of 32 ligands modelled in this entry, 18 are monoatomic - leaving 14 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	1PE	C	405	-	15,15,15	0.62	0	14,14,14	0.99	1 (7%)
3	1PE	E	405	-	15,15,15	0.61	0	14,14,14	0.85	0
3	1PE	C	404	-	15,15,15	0.60	0	14,14,14	0.95	0
3	1PE	F	405	-	15,15,15	0.68	0	14,14,14	0.86	0
3	1PE	D	405	-	15,15,15	0.64	0	14,14,14	0.94	0
3	1PE	A	406	-	15,15,15	0.57	0	14,14,14	0.90	0
3	1PE	A	404	-	15,15,15	0.59	0	14,14,14	0.98	0
3	1PE	A	405	-	15,15,15	0.61	0	14,14,14	0.90	0
3	1PE	C	406	-	15,15,15	0.59	0	14,14,14	0.95	0
3	1PE	D	404	-	15,15,15	0.61	0	14,14,14	0.99	1 (7%)
3	1PE	B	405	-	15,15,15	0.60	0	14,14,14	1.04	2 (14%)
3	1PE	E	404	-	15,15,15	0.61	0	14,14,14	1.01	1 (7%)
3	1PE	B	404	-	15,15,15	0.63	0	14,14,14	0.84	0
3	1PE	F	404	-	15,15,15	0.59	0	14,14,14	0.90	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	1PE	C	405	-	-	9/13/13/13	-
3	1PE	E	405	-	-	9/13/13/13	-
3	1PE	C	404	-	-	9/13/13/13	-
3	1PE	F	405	-	-	8/13/13/13	-
3	1PE	D	405	-	-	5/13/13/13	-
3	1PE	A	406	-	-	8/13/13/13	-
3	1PE	A	404	-	-	5/13/13/13	-
3	1PE	A	405	-	-	8/13/13/13	-
3	1PE	C	406	-	-	9/13/13/13	-
3	1PE	D	404	-	-	6/13/13/13	-
3	1PE	B	405	-	-	7/13/13/13	-
3	1PE	E	404	-	-	9/13/13/13	-
3	1PE	B	404	-	-	7/13/13/13	-
3	1PE	F	404	-	-	10/13/13/13	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	404	1PE	C25-OH5-C14	2.19	122.79	113.29
3	B	405	1PE	C23-OH3-C22	2.19	122.78	113.29
3	E	404	1PE	C25-OH5-C14	2.10	122.37	113.29
3	C	405	1PE	C24-OH4-C13	2.08	122.29	113.29
3	B	405	1PE	C25-OH5-C14	2.02	122.03	113.29

There are no chirality outliers.

All (109) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	404	1PE	C24-C14-OH5-C25
3	B	404	1PE	C16-C26-OH6-C15
3	F	404	1PE	OH6-C15-C25-OH5
3	B	405	1PE	OH6-C15-C25-OH5
3	D	404	1PE	OH5-C14-C24-OH4

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Mol	Chain	Res	Type	Atoms
3	F	405	1PE	OH6-C15-C25-OH5
3	E	405	1PE	OH4-C13-C23-OH3
3	B	405	1PE	OH4-C13-C23-OH3
3	C	404	1PE	OH6-C15-C25-OH5
3	C	406	1PE	OH6-C15-C25-OH5
3	E	404	1PE	OH6-C15-C25-OH5
3	A	406	1PE	OH4-C13-C23-OH3
3	A	406	1PE	OH5-C14-C24-OH4
3	B	404	1PE	OH6-C15-C25-OH5
3	C	405	1PE	OH5-C14-C24-OH4
3	F	404	1PE	C23-C13-OH4-C24
3	C	406	1PE	OH5-C14-C24-OH4
3	C	405	1PE	C25-C15-OH6-C26
3	A	405	1PE	OH5-C14-C24-OH4
3	C	406	1PE	OH4-C13-C23-OH3
3	E	404	1PE	OH4-C13-C23-OH3
3	F	404	1PE	OH7-C16-C26-OH6
3	B	405	1PE	OH2-C12-C22-OH3
3	E	404	1PE	OH7-C16-C26-OH6
3	A	404	1PE	OH7-C16-C26-OH6
3	F	404	1PE	OH4-C13-C23-OH3
3	B	404	1PE	OH4-C13-C23-OH3
3	C	404	1PE	C25-C15-OH6-C26
3	E	404	1PE	OH5-C14-C24-OH4
3	A	406	1PE	OH6-C15-C25-OH5
3	F	404	1PE	OH2-C12-C22-OH3
3	C	404	1PE	OH2-C12-C22-OH3
3	D	404	1PE	OH2-C12-C22-OH3
3	F	404	1PE	OH5-C14-C24-OH4
3	F	405	1PE	OH7-C16-C26-OH6
3	A	406	1PE	C25-C15-OH6-C26
3	C	404	1PE	OH7-C16-C26-OH6
3	B	404	1PE	OH7-C16-C26-OH6
3	A	404	1PE	OH6-C15-C25-OH5
3	D	405	1PE	OH5-C14-C24-OH4
3	B	404	1PE	OH5-C14-C24-OH4
3	E	404	1PE	OH2-C12-C22-OH3
3	C	406	1PE	OH7-C16-C26-OH6
3	C	405	1PE	OH2-C12-C22-OH3
3	E	405	1PE	OH2-C12-C22-OH3
3	C	406	1PE	C25-C15-OH6-C26
3	F	405	1PE	C15-C25-OH5-C14

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Mol	Chain	Res	Type	Atoms
3	B	405	1PE	C16-C26-OH6-C15
3	E	404	1PE	C13-C23-OH3-C22
3	C	405	1PE	C23-C13-OH4-C24
3	C	406	1PE	C14-C24-OH4-C13
3	A	405	1PE	C14-C24-OH4-C13
3	C	406	1PE	C16-C26-OH6-C15
3	A	406	1PE	C24-C14-OH5-C25
3	D	405	1PE	C15-C25-OH5-C14
3	C	404	1PE	C16-C26-OH6-C15
3	C	406	1PE	C15-C25-OH5-C14
3	E	405	1PE	C12-C22-OH3-C23
3	F	405	1PE	C24-C14-OH5-C25
3	F	405	1PE	C16-C26-OH6-C15
3	F	405	1PE	C23-C13-OH4-C24
3	E	405	1PE	OH7-C16-C26-OH6
3	E	405	1PE	C16-C26-OH6-C15
3	D	405	1PE	C23-C13-OH4-C24
3	D	405	1PE	C16-C26-OH6-C15
3	B	404	1PE	C24-C14-OH5-C25
3	E	405	1PE	C23-C13-OH4-C24
3	B	404	1PE	C14-C24-OH4-C13
3	A	404	1PE	OH4-C13-C23-OH3
3	A	406	1PE	C23-C13-OH4-C24
3	D	405	1PE	C12-C22-OH3-C23
3	C	404	1PE	C14-C24-OH4-C13
3	C	405	1PE	C16-C26-OH6-C15
3	A	404	1PE	C16-C26-OH6-C15
3	B	405	1PE	C15-C25-OH5-C14
3	B	405	1PE	OH7-C16-C26-OH6
3	A	406	1PE	OH2-C12-C22-OH3
3	A	405	1PE	C16-C26-OH6-C15
3	A	405	1PE	C24-C14-OH5-C25
3	C	405	1PE	OH4-C13-C23-OH3
3	D	404	1PE	C14-C24-OH4-C13
3	E	405	1PE	C24-C14-OH5-C25
3	F	404	1PE	C25-C15-OH6-C26
3	A	405	1PE	C15-C25-OH5-C14
3	C	404	1PE	OH5-C14-C24-OH4
3	A	405	1PE	OH4-C13-C23-OH3
3	A	405	1PE	OH6-C15-C25-OH5
3	C	405	1PE	OH6-C15-C25-OH5
3	A	405	1PE	C23-C13-OH4-C24

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Mol	Chain	Res	Type	Atoms
3	A	406	1PE	C15-C25-OH5-C14
3	F	405	1PE	C25-C15-OH6-C26
3	B	405	1PE	C12-C22-OH3-C23
3	E	405	1PE	OH6-C15-C25-OH5
3	F	404	1PE	C14-C24-OH4-C13
3	E	404	1PE	C16-C26-OH6-C15
3	E	404	1PE	C24-C14-OH5-C25
3	C	404	1PE	C15-C25-OH5-C14
3	C	404	1PE	C23-C13-OH4-C24
3	C	405	1PE	C24-C14-OH5-C25
3	E	404	1PE	C12-C22-OH3-C23
3	C	405	1PE	C12-C22-OH3-C23
3	D	404	1PE	OH7-C16-C26-OH6
3	F	404	1PE	C12-C22-OH3-C23
3	C	406	1PE	C12-C22-OH3-C23
3	A	404	1PE	OH2-C12-C22-OH3
3	D	404	1PE	C12-C22-OH3-C23
3	F	405	1PE	OH4-C13-C23-OH3
3	E	405	1PE	OH5-C14-C24-OH4
3	D	404	1PE	OH4-C13-C23-OH3

There are no ring outliers.

12 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	405	1PE	3	0
3	E	405	1PE	6	0
3	C	404	1PE	1	0
3	F	405	1PE	4	0
3	A	404	1PE	1	0
3	A	405	1PE	4	0
3	C	406	1PE	1	0
3	D	404	1PE	1	0
3	B	405	1PE	1	0
3	E	404	1PE	4	0
3	B	404	1PE	6	0
3	F	404	1PE	3	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	271/289 (93%)	-0.69	1 (0%)	92	95	5, 10, 21, 28	0
1	B	271/289 (93%)	-0.65	0	100	100	4, 11, 23, 38	0
1	C	271/289 (93%)	-0.65	1 (0%)	92	95	6, 13, 23, 38	0
1	D	271/289 (93%)	-0.68	2 (0%)	87	91	4, 10, 20, 33	0
1	E	271/289 (93%)	-0.59	2 (0%)	87	91	5, 12, 24, 36	0
1	F	271/289 (93%)	-0.62	1 (0%)	92	95	6, 14, 26, 36	0
All	All	1626/1734 (93%)	-0.65	7 (0%)	92	95	4, 12, 24, 38	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	66	GLY	2.5
1	D	114	VAL	2.4
1	D	279	LEU	2.2
1	E	279	LEU	2.2
1	E	337	VAL	2.2
1	A	279	LEU	2.1
1	C	66	GLY	2.1

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

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

6.3 Carbohydrates (i)

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
3	1PE	C	404	16/16	0.82	0.28	33,44,56,62	0
3	1PE	E	404	16/16	0.84	0.19	34,46,53,57	0
3	1PE	D	404	16/16	0.85	0.25	31,40,45,50	0
3	1PE	A	404	16/16	0.86	0.27	28,44,51,54	0
3	1PE	A	405	16/16	0.86	0.24	30,41,49,51	0
3	1PE	C	406	16/16	0.87	0.21	36,45,54,64	0
3	1PE	F	404	16/16	0.88	0.20	33,41,54,62	0
3	1PE	A	406	16/16	0.88	0.27	31,43,54,57	0
3	1PE	E	405	16/16	0.89	0.30	33,43,53,56	0
3	1PE	B	405	16/16	0.89	0.28	37,47,53,57	0
3	1PE	D	405	16/16	0.90	0.31	31,43,50,54	0
3	1PE	F	405	16/16	0.91	0.28	30,35,51,52	0
3	1PE	C	405	16/16	0.92	0.24	26,34,49,52	0
3	1PE	B	404	16/16	0.92	0.24	35,41,51,54	0
2	CA	F	401	1/1	0.98	0.03	12,12,12,12	0
2	CA	D	401	1/1	0.99	0.04	11,11,11,11	0
2	CA	D	402	1/1	0.99	0.06	4,4,4,4	0
2	CA	E	401	1/1	0.99	0.04	11,11,11,11	0
2	CA	E	402	1/1	0.99	0.05	3,3,3,3	0
2	CA	A	401	1/1	0.99	0.04	12,12,12,12	0
2	CA	A	402	1/1	0.99	0.06	5,5,5,5	0
2	CA	F	402	1/1	0.99	0.05	6,6,6,6	0
2	CA	C	401	1/1	0.99	0.03	13,13,13,13	0
2	CA	B	401	1/1	0.99	0.04	13,13,13,13	0
2	CA	B	402	1/1	0.99	0.04	6,6,6,6	0
2	CA	D	403	1/1	1.00	0.03	6,6,6,6	0
2	CA	C	402	1/1	1.00	0.03	8,8,8,8	0
2	CA	B	403	1/1	1.00	0.05	4,4,4,4	0
2	CA	E	403	1/1	1.00	0.03	5,5,5,5	0
2	CA	C	403	1/1	1.00	0.06	5,5,5,5	0
2	CA	F	403	1/1	1.00	0.04	8,8,8,8	0
2	CA	A	403	1/1	1.00	0.04	3,3,3,3	0

6.5 Other polymers (i)

There are no such residues in this entry.