



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:50 PM UTC

PDB ID : 6EWB / pdb_00006ewb
Title : Crystal structure of GII.4 UNSW 2012 P domain in complex with Fab 10E9
Authors : Koromyslova, A.D.; Hansman, G.S.
Deposited on : 2017-11-03
Resolution : 2.78 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

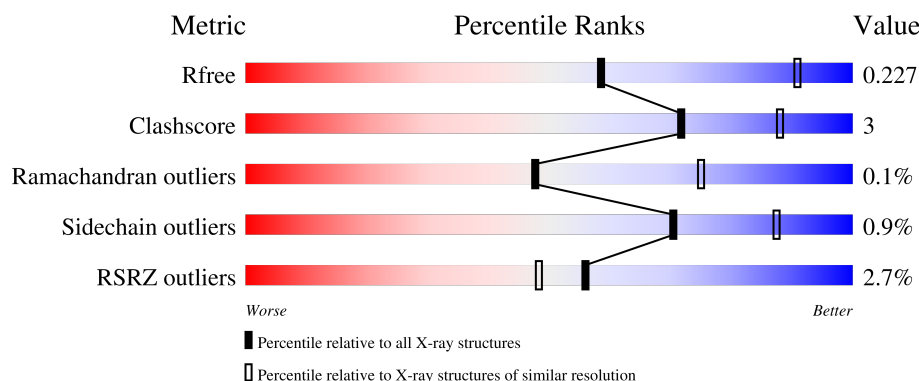
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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}	180053	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	310	10% 92% 6%
1	B	310	89% 10%
1	C	310	93% 7%
1	D	310	93% 7%
2	E	216	10% 84% 6% 10%

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Mol	Chain	Length	Quality of chain
2	G	216	% 88% 9% .
2	H	216	% 87% 10% .
2	J	216	5% 89% 6% .
3	F	214	5% 89% 9% .
3	I	214	2% 87% 11% .
3	K	214	7% 79% 17% .
3	L	214	2% 88% 9% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21709 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 VP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2376	1511	400	455	10			
1	B	306	Total	C	N	O	S	0	0	0
			2384	1512	405	457	10			
1	C	310	Total	C	N	O	S	0	0	0
			2388	1517	402	459	10			
1	D	309	Total	C	N	O	S	0	0	0
			2400	1522	409	459	10			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	221	GLY	-	expression tag	UNP K4LM89
A	222	PRO	-	expression tag	UNP K4LM89
A	223	GLY	-	expression tag	UNP K4LM89
A	224	SER	-	expression tag	UNP K4LM89
B	221	GLY	-	expression tag	UNP K4LM89
B	222	PRO	-	expression tag	UNP K4LM89
B	223	GLY	-	expression tag	UNP K4LM89
B	224	SER	-	expression tag	UNP K4LM89
C	221	GLY	-	expression tag	UNP K4LM89
C	222	PRO	-	expression tag	UNP K4LM89
C	223	GLY	-	expression tag	UNP K4LM89
C	224	SER	-	expression tag	UNP K4LM89
D	221	GLY	-	expression tag	UNP K4LM89
D	222	PRO	-	expression tag	UNP K4LM89
D	223	GLY	-	expression tag	UNP K4LM89
D	224	SER	-	expression tag	UNP K4LM89

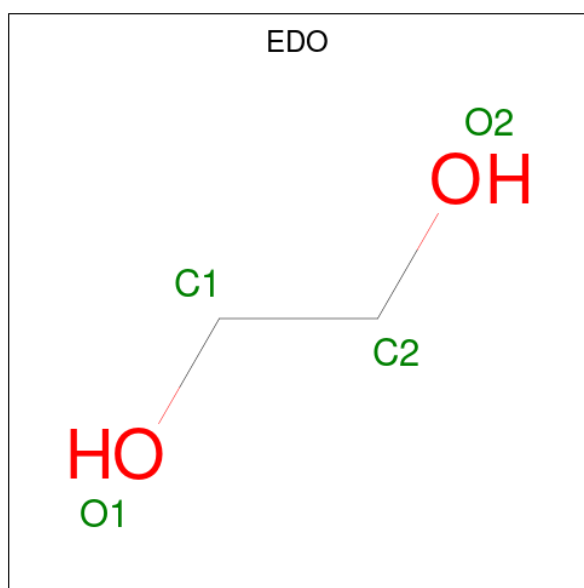
- Molecule 2 is a protein called Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	195	Total	C	N	O	S	0	0	0
			1366	871	222	267	6			
2	G	211	Total	C	N	O	S	0	0	0
			1566	995	257	307	7			
2	H	210	Total	C	N	O	S	0	0	0
			1558	987	255	310	6			
2	J	207	Total	C	N	O	S	0	0	0
			1483	946	242	289	6			

- Molecule 3 is a protein called Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	209	Total	C	N	O	S	0	0	0
			1517	951	252	305	9			
3	I	211	Total	C	N	O	S	0	0	0
			1589	992	261	327	9			
3	K	205	Total	C	N	O	S	0	0	0
			1513	952	248	304	9			
3	L	210	Total	C	N	O	S	0	0	0
			1561	976	258	318	9			

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



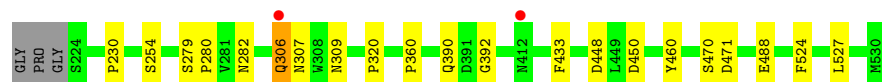
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	L	1	Total	C	O	0	0
			4	2	2		

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.

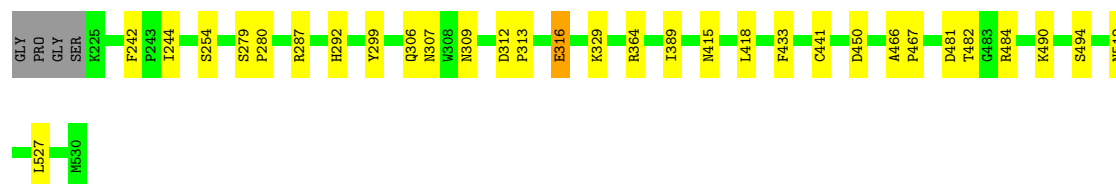
• Molecule 1: VP1

Chain A: 

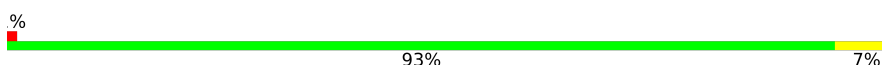


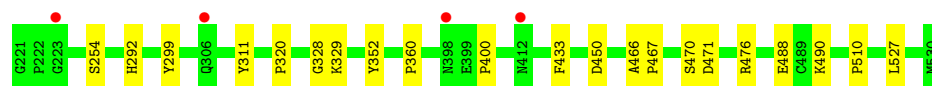
• Molecule 1: VP1

Chain B: 

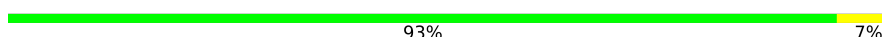


• Molecule 1: VP1

Chain C: 



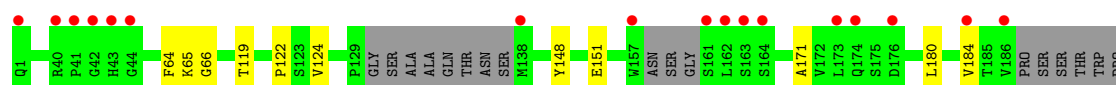
• Molecule 1: VP1

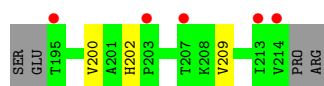
Chain D: 



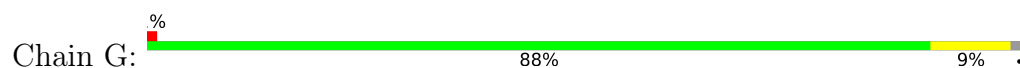
• Molecule 2: Fab heavy chain

Chain E: 

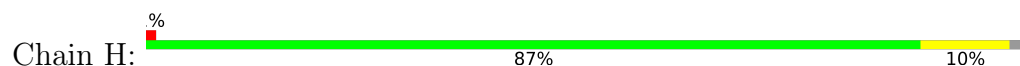




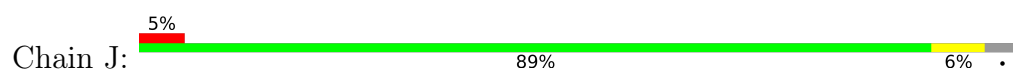
- Molecule 2: Fab heavy chain



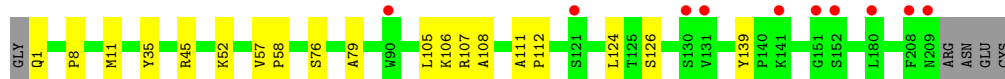
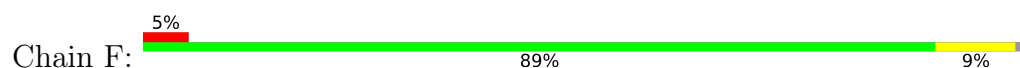
- Molecule 2: Fab heavy chain



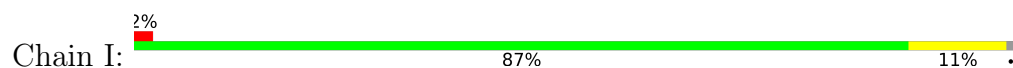
- Molecule 2: Fab heavy chain



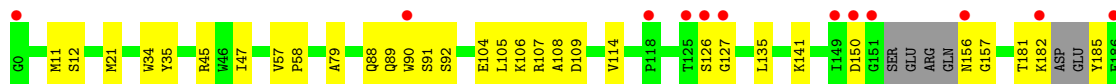
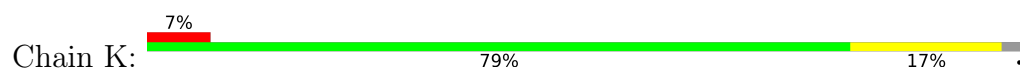
- Molecule 3: Fab light chain

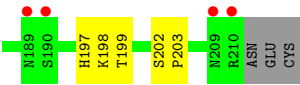


- Molecule 3: Fab light chain

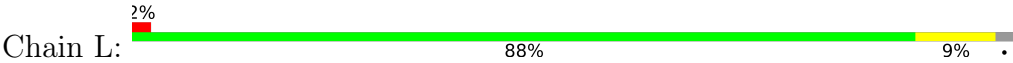


- Molecule 3: Fab light chain





● Molecule 3: Fab light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	106.84Å 111.63Å 288.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.76 – 2.78 48.76 – 2.78	Depositor EDS
% Data completeness (in resolution range)	98.8 (48.76-2.78) 98.9 (48.76-2.78)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.186 , 0.222 0.191 , 0.227	Depositor DCC
R_{free} test set	4336 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21709	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.31% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO

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	0/2446	0.78	0/3349
1	B	0.37	0/2453	0.77	0/3357
1	C	0.36	0/2459	0.77	1/3368 (0.0%)
1	D	0.39	0/2470	0.76	0/3380
2	E	0.36	0/1397	0.80	2/1923 (0.1%)
2	G	0.39	0/1608	0.81	0/2204
2	H	0.43	0/1600	0.87	6/2196 (0.3%)
2	J	0.37	0/1523	0.80	1/2096 (0.0%)
3	F	0.35	0/1555	0.75	0/2128
3	I	0.36	0/1628	0.78	0/2219
3	K	0.36	0/1550	0.76	0/2113
3	L	0.38	0/1600	0.79	2/2185 (0.1%)
All	All	0.38	0/22289	0.78	12/30518 (0.0%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	165	GLY	N-CA-C	-8.08	103.52	115.72
2	E	202	HIS	CA-C-N	6.27	126.47	119.32
2	E	202	HIS	C-N-CA	6.27	126.47	119.32
2	H	52	LEU	CA-C-N	6.17	126.53	119.93
2	H	52	LEU	C-N-CA	6.17	126.53	119.93
2	H	214	VAL	CA-C-N	5.71	126.98	119.84
2	H	214	VAL	C-N-CA	5.71	126.98	119.84
2	H	202	HIS	CA-C-N	5.62	125.73	119.32
2	H	202	HIS	C-N-CA	5.62	125.73	119.32
3	L	57	VAL	CA-C-N	5.59	125.54	119.78
3	L	57	VAL	C-N-CA	5.59	125.54	119.78
1	C	510	PRO	O-C-N	5.02	123.51	121.15

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	2376	0	2258	11	0
1	B	2384	0	2276	16	0
1	C	2388	0	2263	10	0
1	D	2400	0	2284	10	0
2	E	1366	0	1211	6	0
2	G	1566	0	1506	10	0
2	H	1558	0	1486	11	0
2	J	1483	0	1344	9	0
3	F	1517	0	1354	10	0
3	I	1589	0	1480	13	0
3	K	1513	0	1389	22	0
3	L	1561	0	1428	14	0
4	B	4	0	6	0	0
4	L	4	0	6	0	0
All	All	21709	0	20291	139	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 (139) 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)
3:K:89:GLN:HE21	3:K:92:SER:H	1.25	0.83
3:I:89:GLN:HE21	3:I:92:SER:H	1.27	0.80
1:C:329:LYS:HE3	1:C:352:TYR:HB2	1.68	0.74
3:I:89:GLN:NE2	3:I:92:SER:H	1.85	0.74
2:G:154:THR:HB	2:G:201:ALA:HB3	1.72	0.72
3:K:89:GLN:NE2	3:K:92:SER:H	1.89	0.70
3:K:107:ARG:NH1	3:K:108:ALA:O	2.25	0.69
1:B:433:PHE:HB3	1:B:450:ASP:HB3	1.77	0.67
3:I:194:GLU:HG2	3:I:205:VAL:HG22	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:124:LEU:O	3:F:126:SER:HA	1.98	0.64
3:F:79:ALA:HA	3:F:105:LEU:HD21	1.82	0.61
2:E:124:VAL:HG21	2:E:209:VAL:HG21	1.83	0.61
1:A:524:PHE:HB3	2:G:194:GLU:HG3	1.83	0.60
1:A:307:ASN:ND2	1:A:309:ASN:OD1	2.35	0.59
2:H:130:GLY:HA2	2:H:216:ARG:HD2	1.85	0.59
1:B:329:LYS:HD2	1:B:389:ILE:HD11	1.85	0.58
1:D:389:ILE:HG22	1:D:441:CYS:HB2	1.86	0.58
3:F:124:LEU:C	3:F:126:SER:HA	2.30	0.56
3:K:114:VAL:HG22	3:K:135:LEU:HD22	1.87	0.56
2:H:40:ARG:HG3	2:H:92:ALA:HB2	1.89	0.55
2:H:151:GLU:OE1	2:H:152:PRO:HA	2.06	0.55
2:H:67:LYS:NZ	2:H:84:SER:O	2.40	0.55
3:L:89:GLN:HE21	3:L:92:SER:H	1.55	0.54
1:C:490:LYS:HG3	1:C:527:LEU:HD11	1.89	0.54
1:C:433:PHE:HB3	1:C:450:ASP:HB3	1.90	0.54
2:G:199:ASN:ND2	2:G:210:ASP:OD1	2.38	0.54
2:G:138:MET:HE3	2:G:185:THR:HG22	1.88	0.54
3:L:89:GLN:NE2	3:L:92:SER:H	2.06	0.53
1:A:433:PHE:HB3	1:A:450:ASP:HB3	1.91	0.53
3:L:11:MET:HE3	3:L:21:MET:HE2	1.90	0.53
2:H:175:SER:O	2:H:175:SER:OG	2.26	0.52
3:K:181:THR:O	3:K:185:TYR:N	2.43	0.52
3:L:35:TYR:OH	3:L:45:ARG:NH1	2.43	0.52
2:G:151:GLU:OE1	2:G:171:ALA:HB3	2.10	0.51
3:I:107:ARG:NH1	3:I:108:ALA:O	2.44	0.51
1:C:320:PRO:HD2	1:C:360:PRO:HB2	1.93	0.51
3:K:79:ALA:HA	3:K:105:LEU:HD21	1.92	0.51
1:A:448:ASP:OD2	3:F:52:LYS:NZ	2.30	0.51
1:B:415:ASN:HB2	1:B:418:LEU:HD11	1.92	0.51
3:L:120:SER:OG	3:L:122:GLU:OE1	2.29	0.50
2:G:122:PRO:HB3	2:G:148:TYR:HB3	1.92	0.50
1:B:490:LYS:HG3	1:B:527:LEU:HD11	1.94	0.50
2:H:187:PRO:O	2:H:190:THR:OG1	2.30	0.50
3:K:109:ASP:OD2	3:K:198:LYS:HE2	2.12	0.49
1:D:433:PHE:HB3	1:D:450:ASP:HB3	1.94	0.49
1:D:230:PRO:HG3	1:D:460:TYR:CG	2.47	0.49
2:E:171:ALA:HB2	2:E:180:LEU:HB2	1.93	0.49
3:I:12:SER:HB2	3:I:106:LYS:HG2	1.94	0.49
2:J:143:CYS:HB2	2:J:157:TRP:CH2	2.49	0.48
2:J:191:TRP:CE2	2:J:192:PRO:HB3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:287:ARG:HD2	1:D:306:GLN:HA	1.95	0.48
3:K:197:HIS:HB3	3:K:199:THR:HG22	1.95	0.48
1:B:482:THR:OG1	1:B:484:ARG:HG2	2.13	0.48
2:H:191:TRP:CD1	2:H:192:PRO:HA	2.49	0.48
1:C:328:GLY:HA3	1:C:400:PRO:HB3	1.96	0.48
3:L:114:VAL:HG22	3:L:135:LEU:HD22	1.97	0.47
3:K:88:GLN:HE22	3:K:90:TRP:CD1	2.33	0.47
3:K:12:SER:O	3:K:106:LYS:HE3	2.15	0.47
1:D:284:CYS:HA	1:D:324:PRO:HD3	1.97	0.47
3:I:60:ARG:NH2	3:I:80:GLU:OE2	2.36	0.47
3:K:104:GLU:OE2	3:K:141:LYS:NZ	2.40	0.47
3:F:35:TYR:OH	3:F:45:ARG:NH1	2.48	0.46
3:K:156:ASN:HA	3:K:157:GLY:HA3	1.58	0.46
2:J:150:PRO:O	2:J:202:HIS:NE2	2.41	0.46
1:A:488:GLU:HG3	1:A:527:LEU:HD22	1.98	0.46
1:C:470:SER:OG	1:C:471:ASP:N	2.48	0.46
1:A:470:SER:OG	1:A:471:ASP:N	2.48	0.46
1:B:316:GLU:H	1:B:316:GLU:HG2	1.55	0.46
1:A:279:SER:HA	1:A:280:PRO:HD3	1.78	0.46
2:G:40:ARG:HG3	2:G:92:ALA:HB2	1.98	0.46
1:C:292:HIS:HB2	1:C:299:TYR:CE1	2.51	0.45
2:J:191:TRP:CG	2:J:192:PRO:HA	2.52	0.45
2:H:199:ASN:ND2	2:H:210:ASP:OD2	2.42	0.45
3:I:88:GLN:HE22	3:I:90:TRP:CD1	2.35	0.45
3:F:107:ARG:HG2	3:F:108:ALA:N	2.31	0.45
1:B:279:SER:HA	1:B:280:PRO:HD3	1.84	0.44
2:J:151:GLU:HB2	2:J:152:PRO:HA	1.99	0.44
1:B:389:ILE:HG22	1:B:441:CYS:HB2	1.99	0.44
1:B:481:ASP:CG	1:B:512:ASN:HD21	2.26	0.44
1:A:390:GLN:CD	1:A:392:GLY:H	2.25	0.44
1:C:311:TYR:CE1	1:C:320:PRO:HG3	2.52	0.44
2:E:151:GLU:OE2	2:E:171:ALA:HB3	2.17	0.44
3:K:89:GLN:HG2	3:K:91:SER:H	1.81	0.44
3:L:57:VAL:HA	3:L:58:PRO:HD3	1.83	0.44
1:D:488:GLU:HG3	1:D:527:LEU:HD22	1.99	0.44
2:J:149:PHE:HA	2:J:150:PRO:HA	1.71	0.44
1:A:282:ASN:CG	1:A:306:GLN:OE1	2.61	0.44
2:J:122:PRO:HB3	2:J:148:TYR:HB3	2.00	0.44
3:K:35:TYR:OH	3:K:45:ARG:NH1	2.50	0.43
1:B:242:PHE:O	1:B:244:ILE:N	2.49	0.43
2:J:159:SER:O	2:J:159:SER:OG	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:2:VAL:HG21	2:H:98:ARG:NH1	2.32	0.43
3:I:75:ASN:OD1	3:I:76:SER:N	2.51	0.43
3:F:8:PRO:HG2	3:F:11:MET:HB3	2.00	0.43
3:L:12:SER:HA	3:L:104:GLU:O	2.19	0.43
1:B:307:ASN:ND2	1:B:309:ASN:HD22	2.16	0.43
3:K:89:GLN:HG2	3:K:90:TRP:N	2.33	0.43
3:L:89:GLN:HG2	3:L:90:TRP:N	2.33	0.43
1:A:320:PRO:HD2	1:A:360:PRO:HB2	2.01	0.43
1:B:466:ALA:HA	1:B:467:PRO:HD3	1.92	0.43
1:D:279:SER:OG	1:D:281:VAL:HG22	2.19	0.43
1:B:481:ASP:OD1	1:B:512:ASN:ND2	2.48	0.42
1:A:230:PRO:HD3	1:A:460:TYR:CD2	2.54	0.42
1:D:340:THR:O	2:G:103:ASN:HB2	2.20	0.42
2:J:191:TRP:CD2	2:J:192:PRO:HB3	2.54	0.42
1:C:476:ARG:NH1	1:C:488:GLU:OE1	2.53	0.42
1:D:469:GLN:HB2	1:D:520:TRP:CD1	2.54	0.42
2:G:141:LEU:HD11	2:G:191:TRP:CG	2.55	0.42
3:L:188:HIS:O	3:L:210:ARG:NH2	2.52	0.42
3:F:106:LYS:HA	3:F:139:TYR:OH	2.19	0.42
2:H:191:TRP:CG	2:H:192:PRO:HA	2.54	0.42
3:K:202:SER:HA	3:K:203:PRO:HD3	1.91	0.42
3:F:57:VAL:HA	3:F:58:PRO:HD3	1.87	0.42
3:K:11:MET:CE	3:K:21:MET:HB3	2.49	0.42
1:B:292:HIS:HB2	1:B:299:TYR:CE1	2.55	0.42
1:D:225:LYS:HA	1:D:226:PRO:HD3	1.83	0.42
3:K:11:MET:HE1	3:K:21:MET:HB3	2.01	0.42
3:L:11:MET:CE	3:L:21:MET:HE2	2.49	0.42
1:B:312:ASP:HA	1:B:313:PRO:HD3	1.90	0.41
2:E:64:PHE:C	2:E:66:GLY:H	2.28	0.41
2:H:202:HIS:ND1	2:H:205:SER:OG	2.48	0.41
3:K:57:VAL:HA	3:K:58:PRO:HD3	1.87	0.41
3:K:34:TRP:HB2	3:K:47:ILE:HB	2.02	0.41
2:E:122:PRO:HB3	2:E:148:TYR:HB3	2.03	0.41
2:E:200:VAL:HB	2:E:209:VAL:CG2	2.51	0.41
2:G:148:TYR:CE1	2:G:153:VAL:HG13	2.56	0.41
3:I:181:THR:OG1	3:I:184:GLU:HG3	2.20	0.41
3:I:107:ARG:HG2	3:I:108:ALA:N	2.36	0.41
3:I:57:VAL:HA	3:I:58:PRO:HD3	1.94	0.41
3:F:111:ALA:HA	3:F:112:PRO:HD3	1.95	0.41
3:K:182:LYS:HA	3:K:185:TYR:HB3	2.02	0.41
3:L:202:SER:HA	3:L:203:PRO:HD3	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:466:ALA:HA	1:C:467:PRO:HD3	1.92	0.40
3:I:118:PRO:HB3	3:I:208:PHE:CE1	2.55	0.40
3:L:186:GLU:HA	3:L:210:ARG:HH12	1.85	0.40
1:B:287:ARG:HD2	1:B:306:GLN:HA	2.03	0.40
3:L:29:VAL:HG11	3:L:89:GLN:HG3	2.03	0.40
3:I:183:ASP:O	3:I:187:ARG:HG3	2.22	0.40
3:K:11:MET:HE3	3:K:21:MET:HE2	2.03	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	305/310 (98%)	294 (96%)	11 (4%)	0	100	100
1	B	304/310 (98%)	293 (96%)	11 (4%)	0	100	100
1	C	308/310 (99%)	296 (96%)	12 (4%)	0	100	100
1	D	307/310 (99%)	296 (96%)	11 (4%)	0	100	100
2	E	187/216 (87%)	182 (97%)	4 (2%)	1 (0%)	24	52
2	G	207/216 (96%)	202 (98%)	5 (2%)	0	100	100
2	H	206/216 (95%)	202 (98%)	4 (2%)	0	100	100
2	J	201/216 (93%)	199 (99%)	2 (1%)	0	100	100
3	F	207/214 (97%)	199 (96%)	8 (4%)	0	100	100
3	I	209/214 (98%)	201 (96%)	8 (4%)	0	100	100
3	K	199/214 (93%)	191 (96%)	5 (2%)	3 (2%)	8	25
3	L	208/214 (97%)	201 (97%)	7 (3%)	0	100	100
All	All	2848/2960 (96%)	2756 (97%)	88 (3%)	4 (0%)	48	75

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	K	150	ASP
2	E	65	LYS
3	K	127	GLY
3	K	126	SER

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	261/268 (97%)	259 (99%)	2 (1%)	73	89
1	B	264/268 (98%)	260 (98%)	4 (2%)	57	81
1	C	262/268 (98%)	261 (100%)	1 (0%)	84	93
1	D	264/268 (98%)	262 (99%)	2 (1%)	73	89
2	E	129/184 (70%)	127 (98%)	2 (2%)	55	81
2	G	172/184 (94%)	169 (98%)	3 (2%)	53	80
2	H	172/184 (94%)	170 (99%)	2 (1%)	63	84
2	J	148/184 (80%)	147 (99%)	1 (1%)	76	90
3	F	153/188 (81%)	151 (99%)	2 (1%)	61	83
3	I	175/188 (93%)	173 (99%)	2 (1%)	65	85
3	K	159/188 (85%)	159 (100%)	0	100	100
3	L	166/188 (88%)	166 (100%)	0	100	100
All	All	2325/2560 (91%)	2304 (99%)	21 (1%)	70	87

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	254	SER
1	A	306	GLN
1	B	254	SER
1	B	316	GLU
1	B	364	ARG
1	B	494	SER
1	C	254	SER

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Mol	Chain	Res	Type
1	D	254	SER
1	D	486	LEU
2	E	119	THR
2	E	184	VAL
3	F	1	GLN
3	F	76	SER
2	G	89	GLU
2	G	137	SER
2	G	193	SER
2	H	137	SER
2	H	143	CYS
3	I	125	THR
3	I	152	SER
2	J	13	LYS

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

Mol	Chain	Res	Type
1	A	277	GLN
1	A	282	ASN
1	B	277	GLN
1	B	298	ASN
1	B	309	ASN
1	B	501	HIS
1	C	277	GLN
1	C	396	HIS
1	C	463	GLN
1	D	277	GLN
1	D	298	ASN
1	D	302	ASN
1	D	366	GLN
1	D	505	HIS
3	F	136	ASN
3	F	160	ASN
3	I	88	GLN
3	I	189	ASN
3	I	209	ASN
2	J	5	GLN
2	J	59	ASN
3	K	88	GLN
3	K	89	GLN
3	L	1	GLN

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Mol	Chain	Res	Type
3	L	88	GLN
3	L	136	ASN
3	L	160	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	$\# Z > 2$
4	EDO	L	301	-	3,3,3	0.46	0	2,2,2	0.31	0
4	EDO	B	601	-	3,3,3	0.44	0	2,2,2	0.37	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
4	EDO	L	301	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	601	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	307/310 (99%)	-0.39	2 (0%) 84 79	30, 45, 63, 80	0
1	B	306/310 (98%)	-0.38	0 100 100	33, 49, 78, 100	0
1	C	310/310 (100%)	-0.35	4 (1%) 75 69	30, 49, 77, 101	0
1	D	309/310 (99%)	-0.45	1 (0%) 90 88	30, 41, 65, 101	0
2	E	195/216 (90%)	0.72	22 (11%) 10 8	50, 86, 117, 134	0
2	G	211/216 (97%)	-0.23	2 (0%) 81 76	31, 51, 74, 87	0
2	H	210/216 (97%)	-0.26	2 (0%) 79 74	30, 50, 85, 105	0
2	J	207/216 (95%)	0.29	10 (4%) 35 30	40, 71, 117, 147	0
3	F	209/214 (97%)	0.28	10 (4%) 35 30	49, 84, 115, 124	0
3	I	211/214 (98%)	-0.00	4 (1%) 66 59	41, 67, 89, 118	0
3	K	205/214 (95%)	0.26	16 (7%) 19 15	37, 61, 116, 138	0
3	L	210/214 (98%)	-0.15	5 (2%) 59 52	30, 56, 99, 124	0
All	All	2890/2960 (97%)	-0.11	78 (2%) 56 48	30, 55, 104, 147	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	214	VAL	4.9
2	E	161	SER	4.6
3	K	126	SER	4.3
3	K	190	SER	4.2
2	E	157	TRP	4.1
3	K	151	GLY	4.1
2	J	190	THR	3.9
3	K	210	ARG	3.9
2	J	188	SER	3.8
2	J	138	MET	3.7
3	K	150	ASP	3.7

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Mol	Chain	Res	Type	RSRZ
2	E	162	LEU	3.6
3	I	0	GLY	3.4
2	E	176	ASP	3.4
2	G	136	ASN	3.4
3	I	210	ARG	3.2
3	L	156	ASN	3.2
2	E	1	GLN	3.1
3	I	201	THR	3.1
3	K	182	LYS	3.1
2	E	164	SER	3.1
3	K	149	ILE	3.0
3	F	152	SER	3.0
2	E	213	ILE	3.0
2	E	195	THR	2.9
3	K	186	GLU	2.8
2	E	184	VAL	2.8
1	A	306	GLN	2.8
2	E	163	SER	2.8
2	J	176	ASP	2.8
2	E	138	MET	2.6
3	F	90	TRP	2.6
3	F	121	SER	2.6
3	K	90	TRP	2.6
2	E	207	THR	2.6
3	I	90	TRP	2.6
1	A	412	ASN	2.5
2	G	1	GLN	2.5
2	E	43	HIS	2.5
3	F	131	VAL	2.5
3	F	208	PHE	2.5
2	J	194	GLU	2.4
3	K	127	GLY	2.4
2	E	174	GLN	2.4
1	C	412	ASN	2.4
3	L	152	SER	2.4
3	K	0	GLY	2.4
2	J	213	ILE	2.3
2	H	175	SER	2.3
2	J	215	PRO	2.3
3	F	180	LEU	2.3
2	E	41	PRO	2.3
1	C	398	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
3	K	189	ASN	2.3
2	J	137	SER	2.3
3	K	118	PRO	2.3
2	E	173	LEU	2.3
2	E	186	VAL	2.3
3	L	149	ILE	2.2
3	F	209	ASN	2.2
2	J	198	CYS	2.1
2	E	203	PRO	2.1
2	J	54	GLY	2.1
3	L	90	TRP	2.1
3	F	141	LYS	2.1
3	K	156	ASN	2.1
3	F	151	GLY	2.1
3	L	201	THR	2.1
1	D	222	PRO	2.1
3	K	209	ASN	2.1
2	H	130	GLY	2.1
3	F	130	SER	2.0
1	C	223	GLY	2.0
2	E	44	GLY	2.0
3	K	125	THR	2.0
1	C	306	GLN	2.0
2	E	42	GLY	2.0
2	E	40	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
4	EDO	B	601	4/4	0.91	0.16	68,69,69,70	0
4	EDO	L	301	4/4	0.91	0.14	60,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.