



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:55 PM UTC

PDB ID : 4W7A / pdb_00004w7a
Title : Crystal Structure of Full-Length Split GFP Mutant D21H/K26C Disulfide and Metal-Mediated Dimer, P 21 21 21 Space Group, Form 4
Authors : Leibly, D.J.; Waldo, G.S.; Yeates, T.O.
Deposited on : 2014-08-21
Resolution : 3.60 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

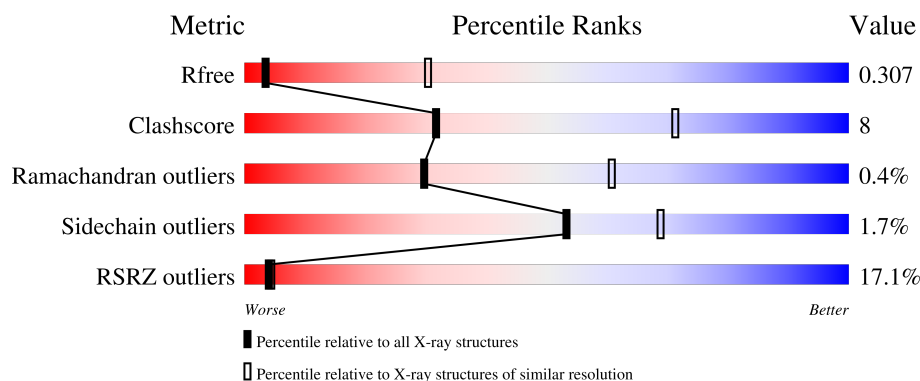
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	229	
1	B	229	
1	C	229	
1	D	229	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	CRO	A	66	-	-	-	X
1	CRO	B	66	-	-	-	X
1	CRO	C	66	-	-	-	X
1	CRO	D	66	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7085 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 fluorescent protein D21H/K26C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	0
			1764	1121	300	338	5			
1	B	223	Total	C	N	O	S	0	0	0
			1786	1137	301	343	5			
1	C	221	Total	C	N	O	S	0	0	0
			1781	1132	302	342	5			
1	D	217	Total	C	N	O	S	0	0	0
			1751	1115	298	334	4			

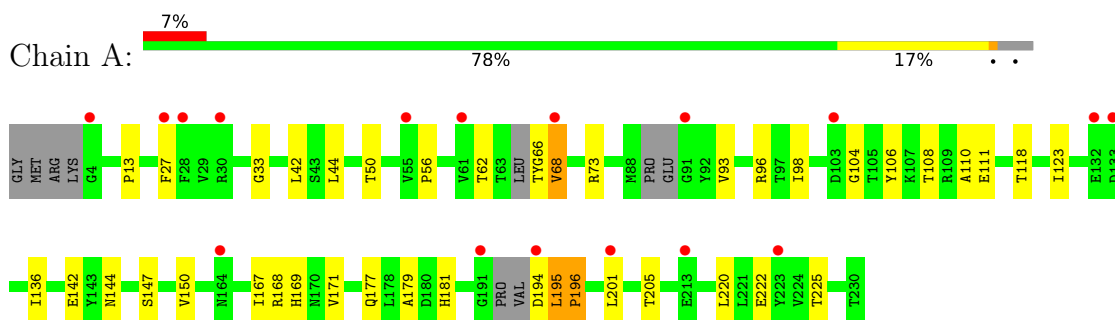
- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cu	0	0
			2	2		
2	C	1	Total	Cu	0	0
			1	1		

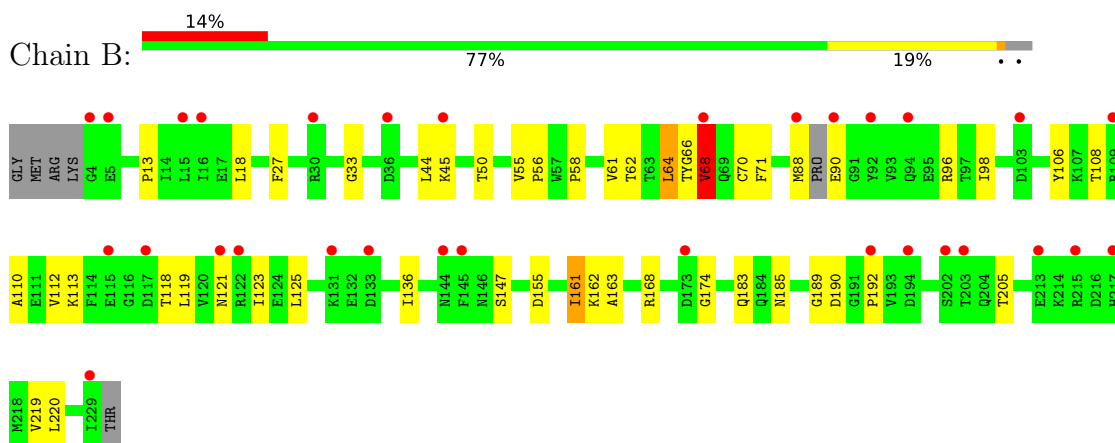
3 Residue-property plots [i](#)

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.

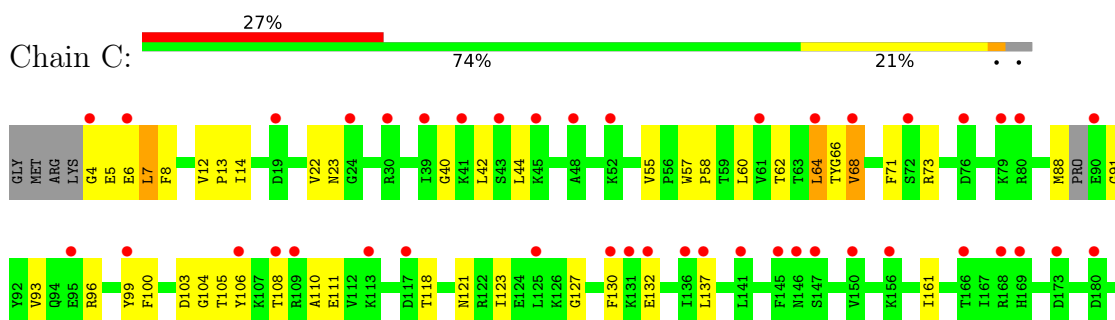
- Molecule 1: fluorescent protein D21H/K26C

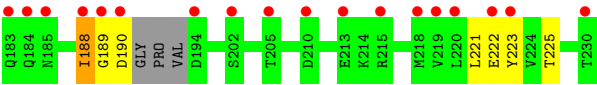


- Molecule 1: fluorescent protein D21H/K26C

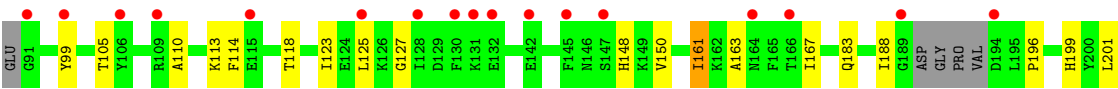
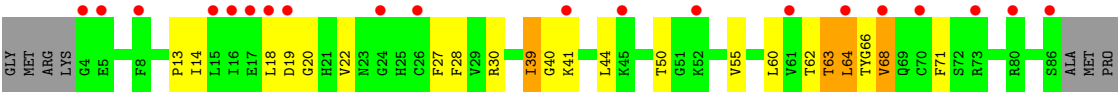
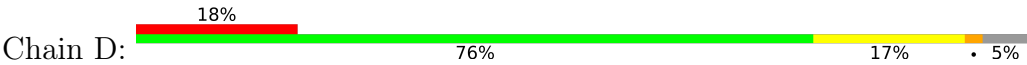


- Molecule 1: fluorescent protein D21H/K26C





● Molecule 1: fluorescent protein D21H/K26C



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	120.83Å 121.33Å 192.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	96.28 – 3.60 96.28 – 3.60	Depositor EDS
% Data completeness (in resolution range)	99.4 (96.28-3.60) 99.4 (96.28-3.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 3.58Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.278 , 0.302 0.285 , 0.307	Depositor DCC
R_{free} test set	3325 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	112.5	Xtriage
Anisotropy	0.432	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 99.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.026 for k,h,-l	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	7085	wwPDB-VP
Average B, all atoms (Å ²)	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% 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: CU, CRO

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	0.32	1/1778 (0.1%)	0.72	1/2400 (0.0%)
1	B	0.35	1/1802 (0.1%)	0.80	4/2436 (0.2%)
1	C	0.33	1/1795 (0.1%)	0.74	3/2423 (0.1%)
1	D	0.37	1/1765 (0.1%)	0.76	3/2383 (0.1%)
All	All	0.34	4/7140 (0.1%)	0.76	11/9642 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	68	VAL	N-CA	-10.80	1.25	1.46
1	A	68	VAL	N-CA	-7.33	1.32	1.46
1	C	68	VAL	N-CA	-6.79	1.33	1.46
1	B	68	VAL	N-CA	-6.13	1.34	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	64	LEU	CA-C-O	-14.30	96.50	120.80
1	D	64	LEU	CA-C-O	-9.36	104.89	120.80
1	C	64	LEU	CA-C-O	-9.18	105.20	120.80
1	B	68	VAL	N-CA-CB	-7.34	99.02	111.50
1	C	7	LEU	N-CA-C	-6.32	105.23	114.39
1	C	6	GLU	N-CA-C	-6.01	101.38	110.70
1	D	63	THR	N-CA-C	5.53	117.11	111.14
1	A	68	VAL	N-CA-CB	-5.40	102.31	111.50
1	D	68	VAL	N-CA-CB	5.36	120.62	111.50
1	B	192	PRO	N-CA-C	-5.18	109.29	114.68
1	B	174	GLY	N-CA-C	-5.18	107.97	115.27

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	1764	0	1702	21	0
1	B	1786	0	1719	31	0
1	C	1781	0	1720	31	0
1	D	1751	0	1694	28	0
2	A	2	0	0	0	0
2	C	1	0	0	0	0
All	All	7085	0	6835	111	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 (111) 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:60:LEU:O	1:C:64:LEU:HG	1.70	0.90
1:C:93:VAL:HG23	1:C:188:ILE:HB	1.65	0.79
1:D:63:THR:HG21	1:D:125:LEU:HD21	1.67	0.76
1:C:5:GLU:C	1:C:7:LEU:H	2.00	0.70
1:D:60:LEU:O	1:D:64:LEU:HG	1.93	0.68
1:C:4:GLY:O	1:C:7:LEU:HB2	1.96	0.66
1:B:98:ILE:HB	1:B:106:TYR:HB2	1.77	0.66
1:A:93:VAL:HG12	1:A:111:GLU:HG2	1.80	0.64
1:B:147:SER:O	1:B:168:ARG:NH2	2.32	0.63
1:D:99:TYR:HA	1:D:105:THR:HG22	1.81	0.63
1:D:20:GLY:HA2	1:D:125:LEU:HB2	1.82	0.62
1:C:188:ILE:O	1:C:188:ILE:HG23	2.00	0.62
1:B:90:GLU:CD	1:B:189:GLY:HA3	2.25	0.61
1:A:68:VAL:O	1:A:68:VAL:HG23	2.01	0.60
1:D:62:THR:HG23	1:D:63:THR:N	2.16	0.60
1:C:13:PRO:HG2	1:C:118:THR:HA	1.83	0.59
1:B:96:ARG:HB2	1:B:108:THR:HB	1.85	0.58
1:C:5:GLU:C	1:C:7:LEU:N	2.62	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:TYR:HA	1:C:105:THR:HG22	1.85	0.57
1:B:68:VAL:HG22	1:B:68:VAL:O	2.03	0.57
1:B:45:LYS:HG2	1:B:219:VAL:HG13	1.85	0.57
1:A:169:HIS:HB2	1:A:177:GLN:HB3	1.87	0.56
1:B:70:CYS:SG	1:B:119:LEU:HD21	2.46	0.56
1:A:98:ILE:HB	1:A:106:TYR:HB2	1.88	0.56
1:C:189:GLY:O	1:C:190:ASP:OD1	2.23	0.56
1:B:205:THR:HG23	1:B:220:LEU:HD11	1.88	0.55
1:D:19:ASP:O	1:D:125:LEU:N	2.33	0.55
1:C:5:GLU:O	1:C:7:LEU:N	2.34	0.55
1:D:18:LEU:HD13	1:D:64:LEU:HD23	1.90	0.54
1:D:110:ALA:HB2	1:D:123:ILE:HG23	1.90	0.54
1:A:27:PHE:HA	1:A:50:THR:HG21	1.88	0.53
1:C:96:ARG:HB2	1:C:108:THR:HB	1.91	0.53
1:B:18:LEU:HD21	1:B:125:LEU:CD2	2.39	0.53
1:C:22:VAL:HG22	1:C:127:GLY:HA3	1.90	0.53
1:B:88:MET:HE1	1:B:113:LYS:HA	1.91	0.52
1:B:163:ALA:HB3	1:B:183:GLN:HG2	1.90	0.52
1:B:33:GLY:HA3	1:B:44:LEU:HD23	1.92	0.52
1:D:39:ILE:HG21	1:D:41:LYS:HE3	1.92	0.52
1:A:62:THR:O	1:A:96:ARG:NH1	2.43	0.51
1:B:161:ILE:HD11	1:B:185:ASN:ND2	2.25	0.51
1:B:161:ILE:HG13	1:B:185:ASN:HB2	1.93	0.51
1:D:13:PRO:HG2	1:D:118:THR:HA	1.93	0.51
1:B:110:ALA:HB2	1:B:123:ILE:HG23	1.93	0.51
1:B:112:VAL:HG22	1:B:121:ASN:HB3	1.92	0.50
1:A:194:ASP:O	1:A:194:ASP:OD1	2.29	0.50
1:A:42:LEU:HB2	1:A:222:GLU:HB2	1.94	0.50
1:C:68:VAL:HG12	1:C:71:PHE:HD2	1.76	0.50
1:A:147:SER:O	1:A:168:ARG:NH2	2.44	0.50
1:D:62:THR:CG2	1:D:63:THR:N	2.75	0.49
1:B:62:THR:O	1:B:96:ARG:NH1	2.44	0.49
1:C:8:PHE:CE1	1:C:88:MET:HG3	2.46	0.49
1:B:205:THR:HG23	1:B:220:LEU:CD1	2.43	0.49
1:D:22:VAL:HG22	1:D:127:GLY:HA3	1.95	0.49
1:B:155:ASP:OD2	1:B:162:LYS:HE3	2.13	0.48
1:C:91:GLY:C	1:C:188:ILE:HG22	2.38	0.48
1:D:27:PHE:HA	1:D:50:THR:HG21	1.94	0.48
1:D:148:HIS:CE1	1:D:167:ILE:HG23	2.49	0.48
1:A:73:ARG:HB3	1:A:225:THR:HG22	1.94	0.48
1:C:93:VAL:HG22	1:C:111:GLU:HG2	1.94	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:PHE:HA	1:B:50:THR:HG21	1.96	0.47
1:A:110:ALA:HB2	1:A:123:ILE:HG23	1.94	0.47
1:B:70:CYS:SG	1:B:119:LEU:CD2	3.02	0.47
1:C:64:LEU:O	1:C:121:ASN:ND2	2.48	0.46
1:A:179:ALA:O	1:A:181:HIS:ND1	2.47	0.46
1:C:221:LEU:HD21	1:C:223:TYR:HE1	1.80	0.46
1:D:63:THR:O	1:D:64:LEU:C	2.59	0.46
1:A:150:VAL:HB	1:A:201:LEU:HB2	1.98	0.46
1:D:148:HIS:HE1	1:D:167:ILE:HG23	1.80	0.46
1:C:8:PHE:HE1	1:C:88:MET:HG3	1.79	0.46
1:D:18:LEU:HD21	1:D:125:LEU:HD12	1.99	0.45
1:D:39:ILE:HB	1:D:41:LYS:HG3	1.97	0.45
1:C:73:ARG:HB3	1:C:225:THR:HG22	1.98	0.45
1:B:71:PHE:CE2	1:B:119:LEU:HD23	2.52	0.44
1:C:14:ILE:HD12	1:C:44:LEU:HD21	1.99	0.44
1:C:12:VAL:HB	1:C:71:PHE:HE1	1.83	0.44
1:A:13:PRO:HG2	1:A:118:THR:HA	2.00	0.44
1:C:42:LEU:HB2	1:C:222:GLU:HB3	2.00	0.44
1:B:18:LEU:HD21	1:B:125:LEU:HD21	1.98	0.43
1:D:41:LYS:HE2	1:D:223:TYR:HE2	1.84	0.43
1:A:62:THR:HG21	1:A:167:ILE:HG13	2.01	0.43
1:A:195:LEU:HA	1:A:196:PRO:HD3	1.69	0.43
1:D:18:LEU:HD13	1:D:64:LEU:CD2	2.48	0.43
1:C:100:PHE:HE1	1:C:106:TYR:HD2	1.66	0.43
1:C:23:ASN:OD1	1:C:130:PHE:HB2	2.17	0.43
1:C:103:ASP:OD1	1:C:104:GLY:N	2.44	0.42
1:C:110:ALA:HB2	1:C:123:ILE:HG23	2.00	0.42
1:C:40:GLY:HA2	1:C:71:PHE:O	2.18	0.42
1:B:56:PRO:HD3	1:B:136:ILE:O	2.19	0.42
1:B:68:VAL:O	1:B:68:VAL:HG13	2.20	0.42
1:C:57:TRP:N	1:C:58:PRO:HD2	2.35	0.42
1:D:161:ILE:HD11	1:D:196:PRO:HG3	2.01	0.42
1:D:163:ALA:HB3	1:D:183:GLN:HB3	2.01	0.42
1:A:96:ARG:HB2	1:A:108:THR:HB	2.00	0.42
1:B:190:ASP:N	1:B:190:ASP:OD1	2.52	0.42
1:D:113:LYS:HE3	1:D:188:ILE:HD13	2.02	0.42
1:D:150:VAL:HB	1:D:201:LEU:HB2	2.01	0.42
1:A:56:PRO:HD3	1:A:136:ILE:O	2.19	0.42
1:A:142:GLU:OE1	1:A:144:ASN:ND2	2.53	0.42
1:B:13:PRO:HG2	1:B:118:THR:HA	2.01	0.42
1:B:58:PRO:HA	1:B:61:VAL:HG23	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:ILE:HD11	1:B:185:ASN:HD22	1.85	0.41
1:C:93:VAL:HG21	1:C:188:ILE:HD13	2.02	0.41
1:D:28:PHE:CE2	1:D:30:ARG:HG3	2.56	0.41
1:B:18:LEU:CD1	1:B:64:LEU:HD21	2.51	0.41
1:D:40:GLY:HA2	1:D:71:PHE:O	2.20	0.41
1:B:96:ARG:HG2	1:B:183:GLN:HB2	2.02	0.41
1:A:33:GLY:HA3	1:A:44:LEU:HD23	2.01	0.41
1:C:132:GLU:HG3	1:C:137:LEU:HD22	2.03	0.41
1:D:14:ILE:HD12	1:D:44:LEU:HD21	2.04	0.41
1:A:205:THR:HG23	1:A:220:LEU:HD11	2.02	0.40
1:D:199:HIS:HB2	1:D:227:ALA:O	2.21	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	211/229 (92%)	202 (96%)	7 (3%)	2 (1%)	14	46
1	B	216/229 (94%)	209 (97%)	7 (3%)	0	100	100
1	C	212/229 (93%)	203 (96%)	8 (4%)	1 (0%)	24	57
1	D	208/229 (91%)	202 (97%)	6 (3%)	0	100	100
All	All	847/916 (92%)	816 (96%)	28 (3%)	3 (0%)	30	61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	GLY
1	C	188	ILE
1	A	196	PRO

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	189/198 (96%)	187 (99%)	2 (1%)	65	74
1	B	192/198 (97%)	189 (98%)	3 (2%)	55	69
1	C	192/198 (97%)	189 (98%)	3 (2%)	55	69
1	D	189/198 (96%)	184 (97%)	5 (3%)	40	62
All	All	762/792 (96%)	749 (98%)	13 (2%)	53	69

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	171	VAL
1	A	195	LEU
1	B	55	VAL
1	B	68	VAL
1	B	161	ILE
1	C	55	VAL
1	C	62	THR
1	C	161	ILE
1	D	39	ILE
1	D	55	VAL
1	D	68	VAL
1	D	114	PHE
1	D	161	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	ASN
1	A	164	ASN
1	A	177	GLN
1	A	184	GLN
1	B	135	ASN
1	C	135	ASN

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Mol	Chain	Res	Type
1	C	159	ASN
1	C	177	GLN
1	D	69	GLN
1	D	94	GLN
1	D	169	HIS
1	D	185	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	CRO	D	66	1	22,23,24	7.09	4 (18%)	30,32,34	3.42	14 (46%)
1	CRO	B	66	1	22,23,24	3.64	6 (27%)	30,32,34	3.46	11 (36%)
1	CRO	A	66	1	22,23,24	3.49	5 (22%)	30,32,34	4.03	9 (30%)
1	CRO	C	66	1	22,23,24	3.73	5 (22%)	30,32,34	2.70	12 (40%)

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
1	CRO	D	66	1	-	4/12/31/32	0/2/2/2
1	CRO	B	66	1	-	3/12/31/32	0/2/2/2
1	CRO	A	66	1	-	4/12/31/32	0/2/2/2
1	CRO	C	66	1	-	5/12/31/32	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	66	CRO	CA3-N3	-32.21	0.87	1.47
1	C	66	CRO	CA3-N3	-15.63	1.18	1.47
1	A	66	CRO	CB2-CA2	14.18	1.48	1.35
1	B	66	CRO	CB2-CA2	14.10	1.48	1.35
1	D	66	CRO	CA1-N1	-6.89	1.28	1.47
1	B	66	CRO	CA3-C3	-5.98	1.28	1.49
1	A	66	CRO	CA3-C3	-5.05	1.31	1.49
1	C	66	CRO	CA1-N1	-4.63	1.34	1.47
1	C	66	CRO	CA3-C3	4.44	1.65	1.49
1	B	66	CRO	CA1-N1	-4.06	1.36	1.47
1	A	66	CRO	C1-N2	3.92	1.37	1.32
1	B	66	CRO	C1-N2	3.88	1.37	1.32
1	B	66	CRO	C2-N3	-2.99	1.33	1.40
1	A	66	CRO	C2-N3	-2.97	1.33	1.40
1	A	66	CRO	O2-C2	2.88	1.28	1.23
1	B	66	CRO	O2-C2	2.88	1.28	1.23
1	D	66	CRO	CA2-N2	2.53	1.44	1.38
1	C	66	CRO	CA2-N2	2.43	1.43	1.38
1	D	66	CRO	C1-N3	2.15	1.40	1.37
1	C	66	CRO	C1-N3	2.08	1.40	1.37

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CRO	C3-CA3-N3	-12.21	84.64	112.43
1	D	66	CRO	O3-C3-CA3	-9.28	83.70	125.47
1	D	66	CRO	C3-CA3-N3	-8.74	92.53	112.43
1	B	66	CRO	O2-C2-CA2	-8.61	125.53	131.02
1	A	66	CRO	O2-C2-CA2	-8.57	125.55	131.02
1	A	66	CRO	CG2-CB2-CA2	-8.30	120.00	129.87
1	B	66	CRO	CG2-CB2-CA2	-8.28	120.02	129.87
1	B	66	CRO	CB2-CA2-C2	7.42	131.36	122.36
1	A	66	CRO	CB2-CA2-C2	7.41	131.34	122.36
1	D	66	CRO	CA3-N3-C2	-6.95	108.40	123.67
1	C	66	CRO	C3-CA3-N3	-6.71	97.15	112.43
1	A	66	CRO	CA2-C2-N3	6.39	108.87	103.50
1	B	66	CRO	CA2-C2-N3	6.39	108.86	103.50
1	A	66	CRO	O3-C3-CA3	-6.02	98.38	125.47
1	D	66	CRO	C1-CA1-N1	5.70	119.96	109.78
1	C	66	CRO	CG2-CB2-CA2	-5.53	123.29	129.87
1	B	66	CRO	C3-CA3-N3	5.51	124.97	112.43
1	D	66	CRO	CA1-C1-N3	-4.97	118.80	124.69

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Mol	Chain	Res	Type	Atoms
1	B	66	CRO	C1-CA1-CB1-OG1
1	B	66	CRO	C1-CA1-CB1-CG1
1	A	66	CRO	N1-CA1-CB1-CG1
1	A	66	CRO	C1-CA1-CB1-OG1
1	A	66	CRO	C1-CA1-CB1-CG1
1	C	66	CRO	N2-C1-CA1-CB1
1	C	66	CRO	N3-C1-CA1-CB1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

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.

