



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 09:24 AM BST

PDB ID : 6VID
Title : The Crystal Structure of Aps Domain-Swapped Trimer Q108K:K40D:T53A:
R58L:Q38F:Q4F Variant of HCRBP2
Authors : Ghanbarpour, A.; Geiger, J.
Deposited on : 2020-01-12
Resolution : 2.89 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	:	2.13
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)	:	2.13

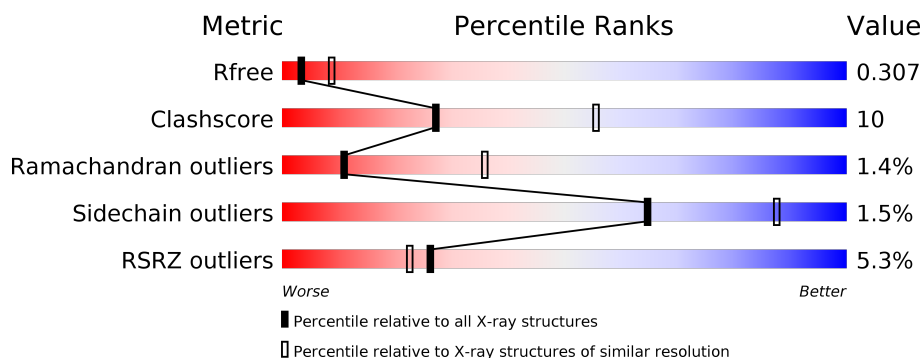
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 $\leq 5\%$. 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	133	<div> <div>73%</div> <div>27%</div> </div>
1	B	133	<div>2%</div> <div>79%</div> <div>20%</div> <div>•</div>
1	C	133	<div>8%</div> <div>65%</div> <div>34%</div> <div>•</div>
1	D	133	<div>82%</div> <div>18%</div>
1	E	133	<div>9%</div> <div>75%</div> <div>24%</div> <div>•</div>
1	F	133	<div>13%</div> <div>74%</div> <div>22%</div> <div>• •</div>

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
3	GOL	D	201	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6167 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 Retinol-binding protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	133	Total	C	N	O	S	0	0	0
			1063	674	176	208	5			
1	B	133	Total	C	N	O	S	0	0	0
			1045	663	173	204	5			
1	C	133	Total	C	N	O	S	0	0	0
			964	610	158	191	5			
1	D	133	Total	C	N	O	S	0	0	0
			1077	684	179	209	5			
1	E	133	Total	C	N	O	S	0	0	0
			1002	638	169	190	5			
1	F	131	Total	C	N	O	S	0	0	0
			954	602	155	192	5			

There are 36 discrepancies between the modelled and reference sequences:

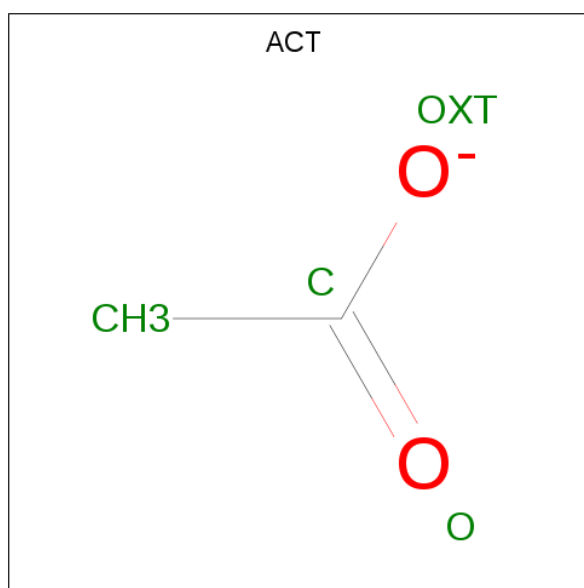
Chain	Residue	Modelled	Actual	Comment	Reference
A	4	PHE	GLN	engineered mutation	UNP P50120
A	38	PHE	GLN	engineered mutation	UNP P50120
A	40	ASP	LYS	engineered mutation	UNP P50120
A	53	ALA	THR	engineered mutation	UNP P50120
A	58	LEU	ARG	engineered mutation	UNP P50120
A	108	LYS	GLN	engineered mutation	UNP P50120
B	4	PHE	GLN	engineered mutation	UNP P50120
B	38	PHE	GLN	engineered mutation	UNP P50120
B	40	ASP	LYS	engineered mutation	UNP P50120
B	53	ALA	THR	engineered mutation	UNP P50120
B	58	LEU	ARG	engineered mutation	UNP P50120
B	108	LYS	GLN	engineered mutation	UNP P50120
C	4	PHE	GLN	engineered mutation	UNP P50120
C	38	PHE	GLN	engineered mutation	UNP P50120
C	40	ASP	LYS	engineered mutation	UNP P50120
C	53	ALA	THR	engineered mutation	UNP P50120
C	58	LEU	ARG	engineered mutation	UNP P50120

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Chain	Residue	Modelled	Actual	Comment	Reference
C	108	LYS	GLN	engineered mutation	UNP P50120
D	4	PHE	GLN	engineered mutation	UNP P50120
D	38	PHE	GLN	engineered mutation	UNP P50120
D	40	ASP	LYS	engineered mutation	UNP P50120
D	53	ALA	THR	engineered mutation	UNP P50120
D	58	LEU	ARG	engineered mutation	UNP P50120
D	108	LYS	GLN	engineered mutation	UNP P50120
E	4	PHE	GLN	engineered mutation	UNP P50120
E	38	PHE	GLN	engineered mutation	UNP P50120
E	40	ASP	LYS	engineered mutation	UNP P50120
E	53	ALA	THR	engineered mutation	UNP P50120
E	58	LEU	ARG	engineered mutation	UNP P50120
E	108	LYS	GLN	engineered mutation	UNP P50120
F	4	PHE	GLN	engineered mutation	UNP P50120
F	38	PHE	GLN	engineered mutation	UNP P50120
F	40	ASP	LYS	engineered mutation	UNP P50120
F	53	ALA	THR	engineered mutation	UNP P50120
F	58	LEU	ARG	engineered mutation	UNP P50120
F	108	LYS	GLN	engineered mutation	UNP P50120

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



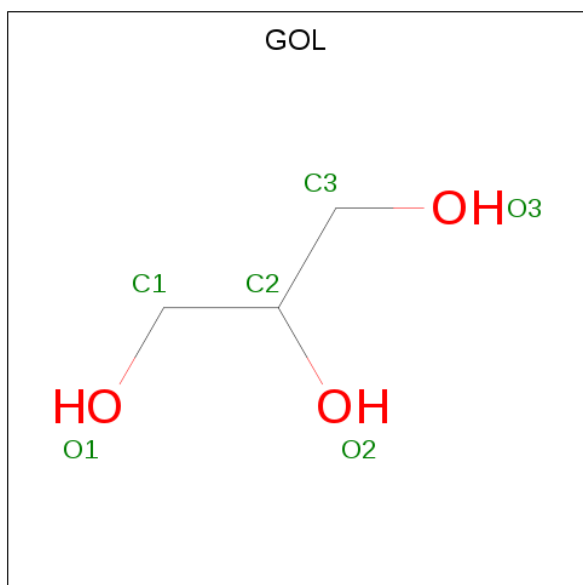
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		
4	D	9	Total	O	0	0
			9	9		

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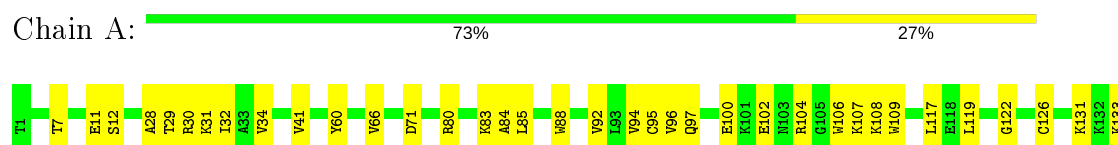
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	2	Total	O	0	0
			2	2		
4	F	1	Total	O	0	0
			1	1		

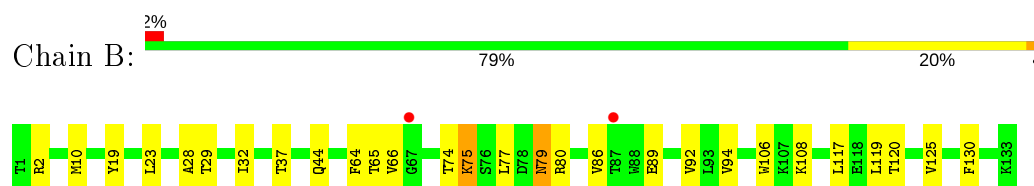
3 Residue-property plots [i](#)

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.

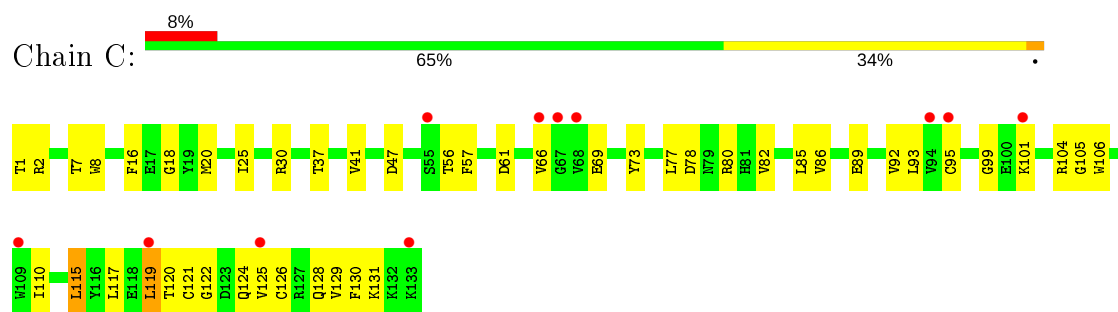
- Molecule 1: Retinol-binding protein 2



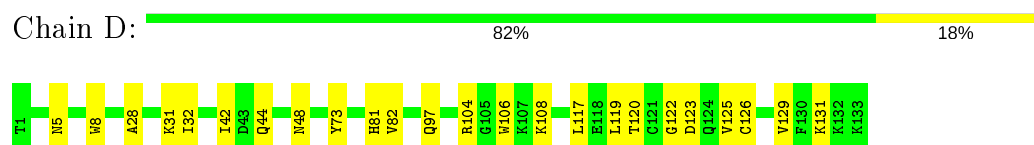
- Molecule 1: Retinol-binding protein 2



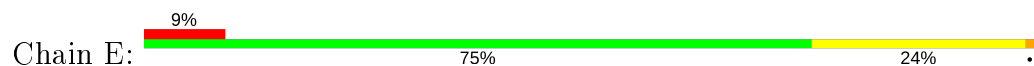
- Molecule 1: Retinol-binding protein 2

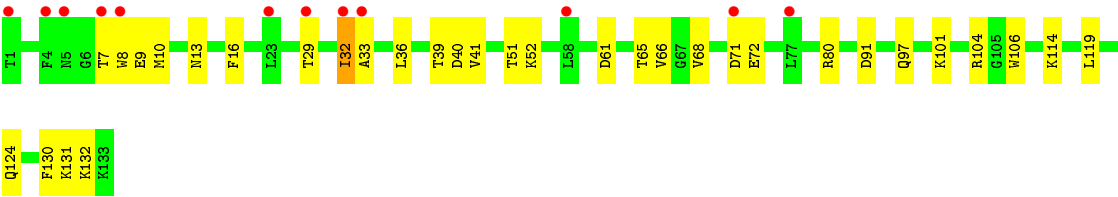


- Molecule 1: Retinol-binding protein 2

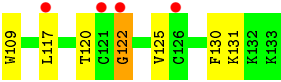
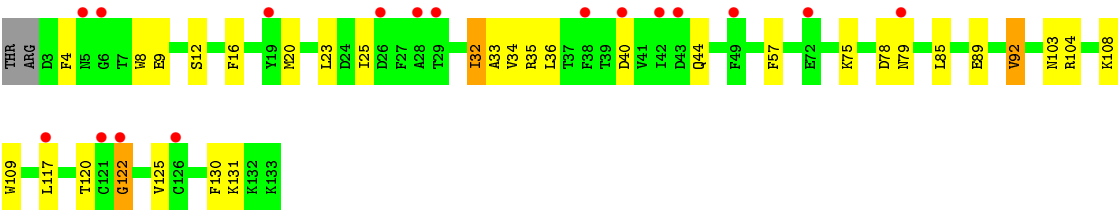


- Molecule 1: Retinol-binding protein 2





• Molecule 1: Retinol-binding protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.05Å 72.68Å 164.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.28 – 2.89 49.28 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.28-2.89) 99.5 (49.28-2.89)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.245 , 0.308 0.245 , 0.307	Depositor DCC
R_{free} test set	1852 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	61.5	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6167	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.29	0/1083	0.53	0/1462
1	B	0.27	0/1064	0.50	0/1440
1	C	0.34	0/983	0.67	2/1344 (0.1%)
1	D	0.26	0/1098	0.47	0/1480
1	E	0.35	0/1021	0.67	1/1387 (0.1%)
1	F	0.33	0/972	0.65	1/1327 (0.1%)
All	All	0.31	0/6221	0.58	4/8440 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	32	ILE	CG1-CB-CG2	-9.88	89.65	111.40
1	C	77	LEU	CB-CG-CD2	7.79	124.25	111.00
1	F	92	VAL	CG1-CB-CG2	5.79	120.17	110.90
1	C	119	LEU	CA-CB-CG	5.01	126.83	115.30

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	1063	0	1004	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1045	0	978	21	0
1	C	964	0	809	31	0
1	D	1077	0	1031	18	0
1	E	1002	0	915	22	0
1	F	954	0	811	22	0
2	A	8	0	6	0	0
2	B	4	0	3	0	0
2	E	4	0	3	0	0
3	A	18	0	24	1	0
3	B	6	0	8	0	0
3	D	6	0	8	0	0
4	A	4	0	0	1	0
4	D	9	0	0	0	0
4	E	2	0	0	0	0
4	F	1	0	0	0	0
All	All	6167	0	5600	119	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 10.

All (119) 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:97:GLN:HB2	1:E:104:ARG:HG2	1.45	0.98
1:B:65:THR:HG22	1:D:48:ASN:HA	1.50	0.90
1:C:115:LEU:HB3	1:C:130:PHE:HB2	1.63	0.80
1:C:104:ARG:HA	1:C:121:CYS:HA	1.62	0.80
1:B:75:LYS:O	1:B:77:LEU:N	2.17	0.77
1:F:20:MET:HA	1:F:23:LEU:HD12	1.71	0.73
1:E:29:THR:HA	1:E:32:ILE:HD12	1.71	0.72
1:B:29:THR:HA	1:B:32:ILE:HD12	1.70	0.72
1:F:89:GLU:O	1:F:92:VAL:HG22	1.92	0.70
1:C:82:VAL:HG12	1:C:99:GLY:HA3	1.74	0.69
1:E:114:LYS:HE3	1:E:131:LYS:HE3	1.75	0.69
1:A:108:LYS:HG3	1:A:117:LEU:HD13	1.75	0.68
1:A:11:GLU:OE2	1:D:131:LYS:HD2	1.96	0.66
1:E:41:VAL:HB	1:E:52:LYS:HB2	1.76	0.65
1:C:110:ILE:HD13	1:F:8:TRP:CH2	2.31	0.64
1:A:131:LYS:HD3	1:A:133:LYS:HE2	1.81	0.63
1:B:75:LYS:C	1:B:77:LEU:H	2.03	0.62
1:C:1:THR:HG22	1:C:2:ARG:H	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:103:ASN:ND2	1:F:122:GLY:O	2.34	0.61
1:D:97:GLN:HB2	1:D:104:ARG:HG2	1.83	0.60
1:E:13:ASN:OD1	1:E:16:PHE:HB2	2.01	0.60
1:B:66:VAL:HA	1:B:86:VAL:HG13	1.85	0.58
1:F:32:ILE:O	1:F:34:VAL:N	2.30	0.58
1:E:80:ARG:HH12	1:E:101:LYS:HE2	1.69	0.57
1:A:60:TYR:HE2	3:A:203:GOL:H2	1.68	0.56
1:A:106:TRP:HB3	1:A:119:LEU:HD23	1.87	0.56
1:B:106:TRP:HB3	1:B:119:LEU:HD23	1.88	0.55
1:C:37:THR:HG22	1:C:57:PHE:O	2.07	0.55
1:A:11:GLU:CD	1:D:131:LYS:HB2	2.27	0.55
1:B:108:LYS:HG3	1:B:117:LEU:HD13	1.88	0.55
1:C:56:THR:OG1	1:C:57:PHE:N	2.38	0.55
1:A:28:ALA:O	1:A:32:ILE:HG13	2.07	0.54
1:C:110:ILE:HD11	1:C:115:LEU:HD12	1.89	0.54
1:F:4:PHE:N	1:F:44:GLN:OE1	2.41	0.54
1:C:120:THR:HG23	1:C:125:VAL:HG22	1.90	0.54
1:A:66:VAL:HG11	1:A:88:TRP:CD1	2.44	0.53
1:D:120:THR:HG23	1:D:125:VAL:HG22	1.91	0.53
1:B:130:PHE:HB3	1:D:8:TRP:HB3	1.90	0.53
1:C:47:ASP:HB3	1:E:66:VAL:HG12	1.90	0.52
1:E:9:GLU:O	1:F:131:LYS:N	2.34	0.52
1:B:2:ARG:O	1:B:44:GLN:NE2	2.43	0.51
1:C:126:CYS:HG	1:F:16:PHE:HD1	1.59	0.51
1:B:74:THR:HG23	1:B:80:ARG:H	1.75	0.51
1:F:32:ILE:O	1:F:34:VAL:HG23	2.11	0.51
1:A:92:VAL:HG22	1:A:109:TRP:HB3	1.92	0.51
1:C:115:LEU:O	1:C:129:VAL:HA	2.11	0.51
1:F:78:ASP:OD2	1:F:104:ARG:NH1	2.44	0.51
1:E:10:MET:HA	1:F:130:PHE:HA	1.93	0.50
1:E:8:TRP:O	1:E:39:THR:HA	2.11	0.50
1:C:95:CYS:O	1:C:105:GLY:HA3	2.11	0.50
1:C:117:LEU:HD22	1:C:128:GLN:NE2	2.26	0.50
1:D:28:ALA:O	1:D:32:ILE:HG13	2.12	0.50
1:A:29:THR:HG23	1:D:31:LYS:HB3	1.93	0.50
1:D:108:LYS:HD3	1:D:117:LEU:HD13	1.94	0.50
1:E:51:THR:O	1:E:52:LYS:HD3	2.12	0.50
1:A:30:ARG:O	1:A:34:VAL:HG13	2.11	0.49
1:C:8:TRP:HB3	1:E:130:PHE:HB3	1.95	0.49
1:B:120:THR:HG22	1:B:125:VAL:HG23	1.93	0.49
1:C:18:GLY:C	1:E:124:GLN:HG2	2.33	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:101:LYS:HA	1:C:101:LYS:HE3	1.93	0.49
1:D:82:VAL:HG21	1:D:97:GLN:HB3	1.95	0.49
1:E:106:TRP:HB3	1:E:119:LEU:HD23	1.94	0.49
1:A:84:ALA:HA	1:A:96:VAL:O	2.13	0.49
1:C:66:VAL:HG21	1:C:93:LEU:HD21	1.94	0.48
1:B:89:GLU:O	1:B:92:VAL:HG22	2.13	0.48
1:C:78:ASP:HA	1:F:25:ILE:HD13	1.96	0.48
1:F:75:LYS:HA	1:F:79:ASN:OD1	2.14	0.47
1:F:34:VAL:O	1:F:36:LEU:N	2.48	0.47
1:A:12:SER:HB2	1:D:129:VAL:HG12	1.97	0.47
1:C:7:THR:HG23	1:C:41:VAL:HG22	1.97	0.47
1:E:101:LYS:HA	1:E:101:LYS:HD3	1.68	0.47
1:A:80:ARG:NH1	1:A:100:GLU:OE2	2.47	0.47
1:E:65:THR:O	1:E:68:VAL:HG22	2.15	0.46
1:F:16:PHE:HE2	1:F:34:VAL:HG12	1.80	0.46
1:A:94:VAL:HG13	1:A:107:LYS:HG2	1.97	0.46
1:C:16:PHE:O	1:C:20:MET:HG3	2.15	0.46
1:C:106:TRP:HB3	1:C:119:LEU:HD23	1.96	0.46
1:F:92:VAL:HG12	1:F:109:TRP:CB	2.45	0.46
1:D:73:TYR:CZ	1:D:81:HIS:HB2	2.51	0.46
1:F:108:LYS:HG3	1:F:117:LEU:HD13	1.98	0.46
1:A:97:GLN:HB2	1:A:104:ARG:HD3	1.98	0.45
1:E:7:THR:HA	1:E:40:ASP:O	2.15	0.45
1:E:72:GLU:OE2	1:E:97:GLN:NE2	2.49	0.45
1:F:120:THR:HG22	1:F:125:VAL:HG23	1.97	0.45
1:D:8:TRP:CD1	1:D:42:ILE:HD12	2.52	0.45
1:E:51:THR:C	1:E:52:LYS:HD3	2.37	0.45
1:C:25:ILE:O	1:C:30:ARG:HD3	2.18	0.44
1:E:33:ALA:HA	1:E:36:LEU:CD1	2.47	0.44
1:A:7:THR:HG23	1:A:41:VAL:HG12	2.00	0.44
1:C:20:MET:HE3	1:C:30:ARG:HG2	2.00	0.44
1:C:69:GLU:HA	1:C:85:LEU:HD12	1.99	0.44
1:C:121:CYS:SG	1:C:122:GLY:N	2.91	0.43
1:A:80:ARG:NH2	1:B:23:LEU:O	2.52	0.43
1:C:131:LYS:N	1:F:9:GLU:O	2.32	0.43
1:A:31:LYS:NZ	4:A:301:HOH:O	2.38	0.43
1:C:89:GLU:O	1:C:92:VAL:HG22	2.19	0.43
1:E:91:ASP:N	1:E:91:ASP:OD1	2.50	0.43
1:A:133:LYS:HA	1:A:133:LYS:HD3	1.83	0.43
1:B:92:VAL:HG23	1:B:94:VAL:HG13	2.01	0.43
1:D:106:TRP:HB3	1:D:119:LEU:HD23	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:LYS:HB3	1:B:29:THR:HG22	2.00	0.42
1:B:28:ALA:O	1:B:32:ILE:HG13	2.19	0.42
1:F:92:VAL:HG12	1:F:109:TRP:HB3	2.01	0.42
1:F:78:ASP:OD1	1:F:78:ASP:N	2.43	0.42
1:C:130:PHE:HB3	1:F:8:TRP:HB3	2.02	0.42
1:B:74:THR:HG22	1:B:80:ARG:O	2.20	0.42
1:A:102:GLU:HG2	1:A:102:GLU:H	1.65	0.42
1:A:126:CYS:HB2	1:B:19:TYR:HB2	2.02	0.41
1:B:75:LYS:HG2	1:B:79:ASN:HD21	1.85	0.41
1:B:64:PHE:CD1	1:B:86:VAL:HG11	2.54	0.41
1:C:80:ARG:NH1	1:C:101:LYS:HD3	2.35	0.41
1:A:11:GLU:OE2	1:D:131:LYS:HB2	2.19	0.41
1:D:5:ASN:OD1	1:D:44:GLN:N	2.38	0.41
1:D:73:TYR:CE2	1:D:81:HIS:HB2	2.55	0.41
1:A:71:ASP:OD1	1:A:83:LYS:HD3	2.20	0.41
1:D:123:ASP:N	1:D:123:ASP:OD1	2.46	0.41
1:A:85:LEU:O	1:A:95:CYS:HA	2.20	0.41
1:C:8:TRP:CH2	1:E:132:LYS:HD3	2.56	0.41
1:B:10:MET:HB3	1:B:37:THR:HG22	2.02	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	131/133 (98%)	127 (97%)	3 (2%)	1 (1%)	19	51
1	B	131/133 (98%)	122 (93%)	8 (6%)	1 (1%)	19	51
1	C	131/133 (98%)	118 (90%)	11 (8%)	2 (2%)	10	34
1	D	131/133 (98%)	128 (98%)	2 (2%)	1 (1%)	19	51
1	E	131/133 (98%)	124 (95%)	6 (5%)	1 (1%)	19	51

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	129/133 (97%)	118 (92%)	6 (5%)	5 (4%)	3	12
All	All	784/798 (98%)	737 (94%)	36 (5%)	11 (1%)	11	36

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	124	GLN
1	F	33	ALA
1	A	122	GLY
1	B	75	LYS
1	C	86	VAL
1	E	61	ASP
1	F	57	PHE
1	F	122	GLY
1	F	35	ARG
1	F	32	ILE
1	D	122	GLY

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	112/119 (94%)	112 (100%)	0	100	100
1	B	108/119 (91%)	107 (99%)	1 (1%)	78	93
1	C	87/119 (73%)	84 (97%)	3 (3%)	37	71
1	D	115/119 (97%)	114 (99%)	1 (1%)	78	93
1	E	97/119 (82%)	96 (99%)	1 (1%)	76	92
1	F	89/119 (75%)	86 (97%)	3 (3%)	37	71
All	All	608/714 (85%)	599 (98%)	9 (2%)	65	87

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	79	ASN
1	C	61	ASP
1	C	73	TYR
1	C	115	LEU
1	D	126	CYS
1	E	71	ASP
1	F	12	SER
1	F	40	ASP
1	F	85	LEU

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

Mol	Chain	Res	Type
1	B	79	ASN
1	C	128	GLN
1	E	79	ASN

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](#)

9 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
3	GOL	B	202	-	5,5,5	0.90	0	5,5,5	0.99	0
3	GOL	A	203	-	5,5,5	0.91	0	5,5,5	1.00	0
3	GOL	A	204	-	5,5,5	0.90	0	5,5,5	1.00	0
2	ACT	A	202	-	1,3,3	6.41	1 (100%)	0,3,3	0.00	-
2	ACT	E	201	-	1,3,3	6.49	1 (100%)	0,3,3	0.00	-
2	ACT	A	201	-	1,3,3	6.37	1 (100%)	0,3,3	0.00	-
2	ACT	B	201	-	1,3,3	6.45	1 (100%)	0,3,3	0.00	-
3	GOL	A	205	-	5,5,5	0.90	0	5,5,5	1.00	0
3	GOL	D	201	-	5,5,5	0.91	0	5,5,5	0.99	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	GOL	B	202	-	-	4/4/4/4	-
3	GOL	A	204	-	-	0/4/4/4	-
3	GOL	A	205	-	-	3/4/4/4	-
3	GOL	A	203	-	-	0/4/4/4	-
3	GOL	D	201	-	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	201	ACT	CH3-C	6.49	1.57	1.48
2	B	201	ACT	CH3-C	6.45	1.57	1.48
2	A	202	ACT	CH3-C	6.41	1.56	1.48
2	A	201	ACT	CH3-C	6.37	1.56	1.48

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	202	GOL	O1-C1-C2-C3
3	B	202	GOL	C1-C2-C3-O3
3	D	201	GOL	O1-C1-C2-C3
3	B	202	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	B	202	GOL	O1-C1-C2-O2
3	A	205	GOL	O2-C2-C3-O3
3	A	205	GOL	C1-C2-C3-O3
3	D	201	GOL	O1-C1-C2-O2
3	A	205	GOL	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	203	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	133/133 (100%)	-0.01	0 100 100	28, 43, 60, 68	0
1	B	133/133 (100%)	0.09	2 (1%) 73 73	29, 50, 69, 78	0
1	C	133/133 (100%)	0.63	11 (8%) 11 8	51, 77, 98, 109	0
1	D	133/133 (100%)	-0.01	0 100 100	31, 44, 66, 72	0
1	E	133/133 (100%)	0.47	12 (9%) 9 7	45, 73, 92, 105	0
1	F	131/133 (98%)	0.78	17 (12%) 3 2	62, 79, 103, 112	0
All	All	796/798 (99%)	0.32	42 (5%) 26 22	28, 61, 94, 112	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	133	LYS	5.9
1	E	1	THR	5.5
1	E	4	PHE	4.7
1	F	26	ASP	4.7
1	C	67	GLY	4.5
1	F	49	PHE	4.4
1	E	29	THR	4.4
1	F	6	GLY	3.6
1	C	94	VAL	3.0
1	C	119	LEU	3.0
1	B	67	GLY	3.0
1	C	55	SER	3.0
1	F	19	TYR	2.9
1	F	38	PHE	2.9
1	F	28	ALA	2.9
1	F	29	THR	2.8
1	E	23	LEU	2.6
1	F	5	ASN	2.6
1	F	40	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	125	VAL	2.5
1	F	79	ASN	2.5
1	C	66	VAL	2.5
1	C	101	LYS	2.5
1	F	43	ASP	2.5
1	C	95	CYS	2.4
1	E	32	ILE	2.4
1	F	42	ILE	2.4
1	F	121	CYS	2.4
1	E	33	ALA	2.4
1	F	122	GLY	2.3
1	E	71	ASP	2.3
1	F	72	GLU	2.3
1	B	87	THR	2.2
1	E	8	TRP	2.2
1	F	126	CYS	2.2
1	E	7	THR	2.1
1	C	109	TRP	2.1
1	E	58	LEU	2.1
1	E	5	ASN	2.1
1	C	68	VAL	2.0
1	E	77	LEU	2.0
1	F	117	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	D	201	6/6	0.62	0.47	56,65,68,68	0
2	ACT	E	201	4/4	0.71	0.39	62,64,78,79	0
3	GOL	A	203	6/6	0.73	0.38	55,62,68,69	0
3	GOL	B	202	6/6	0.78	0.30	50,52,55,63	0
2	ACT	A	202	4/4	0.80	0.39	40,45,52,59	0
2	ACT	B	201	4/4	0.81	0.22	63,64,68,78	0
3	GOL	A	204	6/6	0.85	0.21	52,57,70,73	0
3	GOL	A	205	6/6	0.88	0.27	35,44,49,51	0
2	ACT	A	201	4/4	0.95	0.18	41,53,56,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.