Full wwPDB X-ray Structure Validation Report (i)

Oct 26, 2023 – 04:15 PM EDT

PDB ID	:	3JUT
Title	:	Acidic Fibroblast Growth Factor (FGF-1) complexed with gentisic acid
Authors	:	Fernandez, I.S.; Gimenez-Gallego, G.; Romero, A.
Deposited on	:	2009-09-15
Resolution	:	2.25 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:

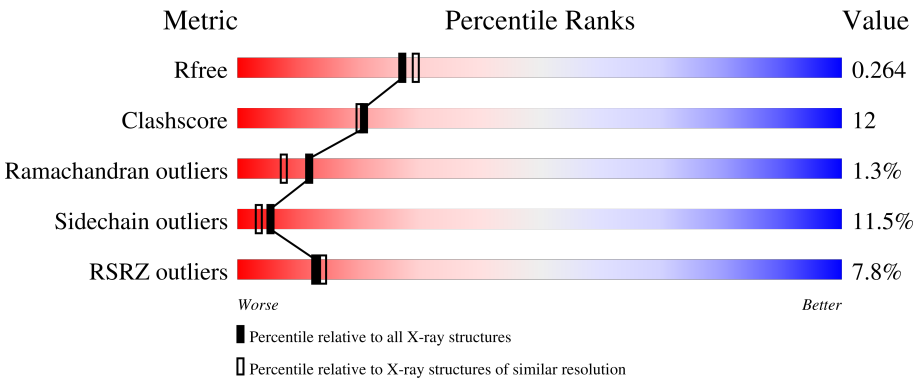
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	130	9% 69% 23% 6% •
1	B	130	5% 75% 20% 5% •
1	C	130	5% 75% 21% •
1	D	130	5% 76% 19% • •
1	E	130	11% 68% 25% 6% •

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Mol	Chain	Length	Quality of chain
1	F	130	11% 52% 36% 11%

2 Entry composition (i)

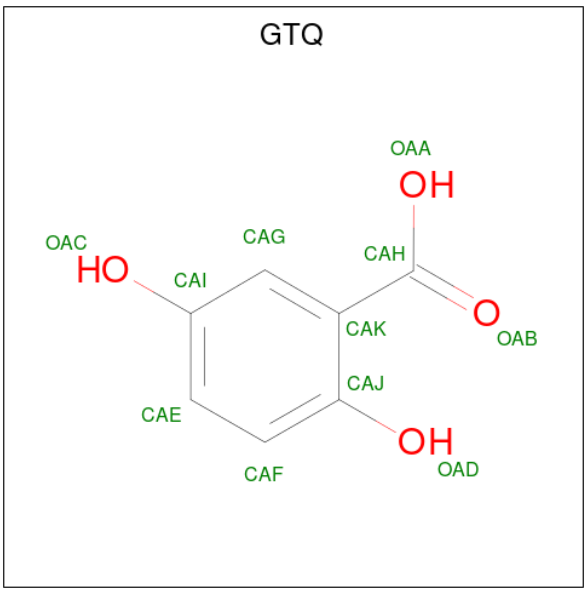
There are 3 unique types of molecules in this entry. The entry contains 6272 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Heparin-binding growth factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	130	1038	655	182	197	4	0	0	0
1	B	130	1037	655	182	196	4	0	0	0
1	C	130	1038	655	182	197	4	0	0	0
1	D	130	1038	655	182	197	4	0	0	0
1	E	130	1038	655	182	197	4	0	0	0
1	F	130	1038	655	182	197	4	0	0	0

- Molecule 2 is 2,5-dihydroxybenzoic acid (three-letter code: GTQ) (formula: $C_7H_6O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	11	7	4	0	0
			Total	C	O		
2	B	1	11	7	4	0	0

- Molecule 3 is water.

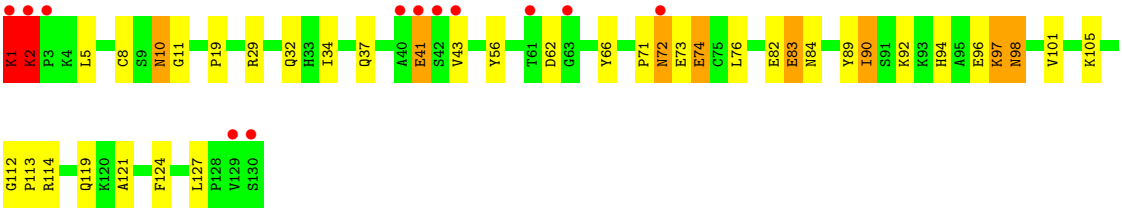
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	6	6	6	0	0
			Total	O		
3	B	5	5	5	0	0
			Total	O		
3	C	4	4	4	0	0
			Total	O		
3	D	6	6	6	0	0
			Total	O		
3	E	2	2	2	0	0

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.

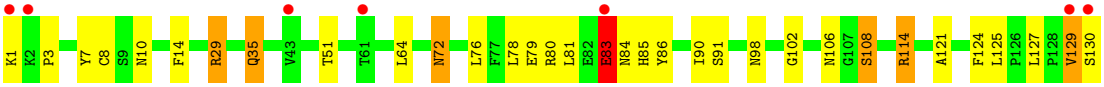
• Molecule 1: Heparin-binding growth factor 1

Chain A:



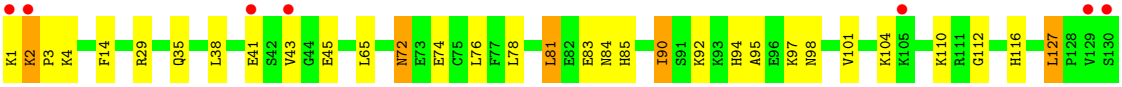
• Molecule 1: Heparin-binding growth factor 1

Chain B:



• Molecule 1: Heparin-binding growth factor 1

Chain C:

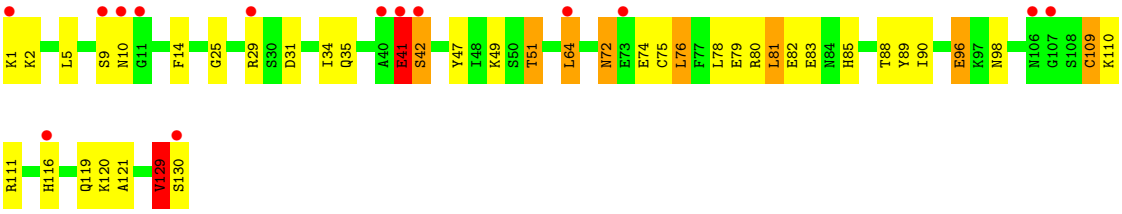


• Molecule 1: Heparin-binding growth factor 1

Chain D:

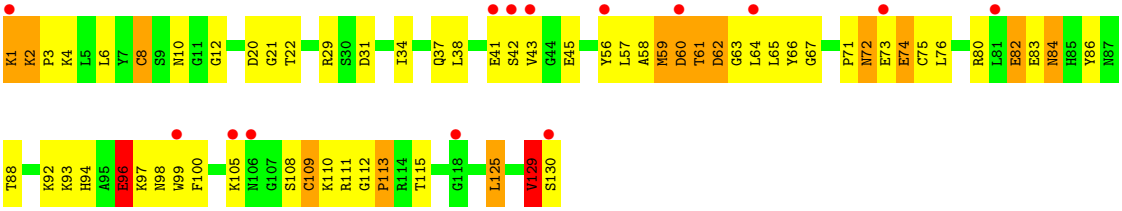
• Molecule 1: Heparin-binding growth factor 1

Chain E:



- Molecule 1: Heparin-binding growth factor 1

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	97.66Å 47.69Å 98.41Å 90.00° 106.45° 90.00°	Depositor
Resolution (Å)	37.01 – 2.25 37.01 – 2.25	Depositor EDS
% Data completeness (in resolution range)	96.2 (37.01-2.25) 96.2 (37.01-2.25)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.12 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.216 , 0.274 0.212 , 0.264	Depositor DCC
R_{free} test set	2034 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	39.1	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 63.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.014 for l,-k,h	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6272	wwPDB-VP
Average B, all atoms (Å^2)	38.0	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 56.30 % 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 2.7915e-05. 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: GTQ

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	1.99	19/1061 (1.8%)	1.17	8/1429 (0.6%)
1	B	1.05	0/1060	0.96	1/1429 (0.1%)
1	C	0.92	0/1061	0.97	2/1429 (0.1%)
1	D	1.02	0/1061	0.94	2/1429 (0.1%)
1	E	0.97	1/1061 (0.1%)	0.94	0/1429
1	F	1.61	23/1061 (2.2%)	1.15	5/1429 (0.3%)
All	All	1.32	43/6365 (0.7%)	1.03	18/8574 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	2
1	F	0	1
All	All	0	3

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	96	GLU	CD-OE2	28.11	1.56	1.25
1	A	83	GLU	CD-OE2	26.08	1.54	1.25
1	A	83	GLU	CD-OE1	21.10	1.48	1.25
1	F	60	ASP	CG-OD2	18.11	1.67	1.25
1	F	74	GLU	CD-OE2	12.33	1.39	1.25
1	A	96	GLU	CG-CD	10.81	1.68	1.51
1	F	21	GLY	C-O	10.42	1.40	1.23
1	F	66	TYR	CE2-CZ	9.61	1.51	1.38
1	A	96	GLU	CD-OE1	9.60	1.36	1.25
1	F	66	TYR	CG-CD1	9.59	1.51	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	73	GLU	CG-CD	9.52	1.66	1.51
1	A	97	LYS	CD-CE	9.29	1.74	1.51
1	A	82	GLU	C-O	8.90	1.40	1.23
1	A	41	GLU	CD-OE2	8.88	1.35	1.25
1	A	2	LYS	CE-NZ	8.74	1.70	1.49
1	A	83	GLU	CG-CD	8.63	1.64	1.51
1	A	41	GLU	CD-OE1	8.48	1.34	1.25
1	A	11	GLY	N-CA	8.15	1.58	1.46
1	F	66	TYR	CE1-CZ	7.92	1.48	1.38
1	F	59	MET	CB-CG	7.72	1.76	1.51
1	F	66	TYR	CG-CD2	7.67	1.49	1.39
1	A	73	GLU	CD-OE2	7.42	1.33	1.25
1	F	96	GLU	C-O	7.24	1.37	1.23
1	F	82	GLU	CG-CD	6.80	1.62	1.51
1	A	1	LYS	CD-CE	6.31	1.67	1.51
1	F	112	GLY	CA-C	6.29	1.61	1.51
1	F	73	GLU	CD-OE1	6.11	1.32	1.25
1	F	96	GLU	CG-CD	6.10	1.61	1.51
1	F	59	MET	C-O	5.99	1.34	1.23
1	F	73	GLU	CD-OE2	5.88	1.32	1.25
1	A	97	LYS	CE-NZ	5.86	1.63	1.49
1	E	75	CYS	CB-SG	-5.70	1.72	1.81
1	F	113	PRO	C-N	5.63	1.47	1.34
1	F	8	CYS	CB-SG	-5.51	1.72	1.81
1	F	92	LYS	CE-NZ	5.42	1.62	1.49
1	F	113	PRO	C-O	5.38	1.34	1.23
1	A	74	GLU	CD-OE2	5.30	1.31	1.25
1	F	62	ASP	CG-OD2	5.30	1.37	1.25
1	A	98	ASN	CB-CG	5.25	1.63	1.51
1	F	93	LYS	CD-CE	5.08	1.64	1.51
1	F	82	GLU	CB-CG	5.08	1.61	1.52
1	F	84	ASN	CB-CG	5.07	1.62	1.51
1	A	94	HIS	C-O	5.05	1.32	1.23

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	F	62	ASP	CB-CG-OD1	15.70	132.43	118.30
1	A	96	GLU	CG-CD-OE2	-12.45	93.41	118.30
1	A	10	ASN	C-N-CA	-10.16	100.97	122.30
1	A	96	GLU	OE1-CD-OE2	7.99	132.88	123.30
1	F	62	ASP	OD1-CG-OD2	-7.45	109.14	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	96	GLU	CG-CD-OE1	7.38	133.06	118.30
1	F	92	LYS	CD-CE-NZ	-7.29	94.92	111.70
1	F	60	ASP	CB-CG-OD2	-6.57	112.39	118.30
1	D	127	LEU	CB-CG-CD1	6.24	121.60	111.00
1	F	61	THR	CA-CB-CG2	-6.04	103.94	112.40
1	C	127	LEU	CA-CB-CG	5.99	129.09	115.30
1	A	11	GLY	N-CA-C	5.90	127.86	113.10
1	A	62	ASP	CB-CG-OD1	5.40	123.16	118.30
1	A	10	ASN	O-C-N	-5.32	114.15	123.20
1	D	29	ARG	NE-CZ-NH2	-5.29	117.66	120.30
1	B	114	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	C	127	LEU	CB-CG-CD1	5.23	119.89	111.00
1	A	29	ARG	CG-CD-NE	5.17	122.66	111.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	129	VAL	Peptide
1	E	41	GLU	Peptide
1	F	129	VAL	Peptide

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	1038	0	1028	25	0
1	B	1037	0	1028	25	0
1	C	1038	0	1028	14	0
1	D	1038	0	1028	20	0
1	E	1038	0	1028	28	0
1	F	1038	0	1028	46	0
2	A	11	0	3	3	0
2	B	11	0	3	1	0
3	A	6	0	0	0	0
3	B	5	0	0	0	0
3	C	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	6	0	0	0	0
3	E	2	0	0	0	0
All	All	6272	0	6174	151	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 12.

All (151) 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:F:59:MET:CB	1:F:59:MET:CG	1.76	1.61
1:A:97:LYS:CD	1:A:97:LYS:CE	1.74	1.59
1:A:2:LYS:CE	1:A:2:LYS:NZ	1.70	1.49
1:F:60:ASP:CG	1:F:60:ASP:OD2	1.67	1.33
1:B:83:GLU:HA	1:B:83:GLU:OE1	1.64	0.94
1:F:96:GLU:H	1:F:96:GLU:CD	1.74	0.90
1:A:119:GLN:HG2	2:A:131:GTQ:HAF	1.51	0.90
1:F:29:ARG:HE	1:F:130:SER:HB2	1.40	0.86
1:D:90:ILE:HD11	1:D:98:ASN:HA	1.62	0.81
1:B:29:ARG:HH11	1:B:29:ARG:CB	1.95	0.80
1:B:29:ARG:HH11	1:B:29:ARG:HB2	1.47	0.79
1:C:72:ASN:HD22	1:C:74:GLU:H	1.30	0.78
1:A:97:LYS:CE	1:A:97:LYS:CG	2.62	0.77
1:F:8:CYS:O	1:F:12:GLY:HA2	1.85	0.75
1:D:72:ASN:HD22	1:D:74:GLU:H	1.37	0.72
1:B:10:ASN:OD1	2:B:131:GTQ:HAE	1.88	0.72
1:A:90:ILE:HD12	1:A:98:ASN:HA	1.71	0.72
1:F:59:MET:CG	1:F:59:MET:CA	2.66	0.71
1:B:129:VAL:O	1:B:130:SER:HB2	1.91	0.71
1:B:83:GLU:OE1	1:B:83:GLU:CA	2.38	0.70
1:D:3:PRO:HG2	1:D:129:VAL:HG11	1.74	0.70
1:E:29:ARG:HH11	1:E:29:ARG:HB3	1.56	0.70
1:B:79:GLU:O	1:B:80:ARG:NH1	2.27	0.67
1:E:41:GLU:CG	1:E:42:SER:HB2	2.24	0.67
1:F:72:ASN:HD22	1:F:74:GLU:H	1.42	0.66
1:E:81:LEU:HD22	1:E:85:HIS:HA	1.77	0.66
1:C:72:ASN:ND2	1:C:74:GLU:H	1.94	0.65
1:A:97:LYS:CD	1:A:97:LYS:NZ	2.60	0.65
1:D:78:LEU:HB2	1:D:90:ILE:HG22	1.77	0.65
1:F:63:GLY:HA3	1:F:113:PRO:HD3	1.78	0.64
1:B:129:VAL:O	1:B:129:VAL:HG22	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:72:ASN:ND2	1:F:74:GLU:H	1.96	0.63
1:F:72:ASN:HD22	1:F:72:ASN:C	2.04	0.61
1:F:111:ARG:HB3	1:F:113:PRO:HD2	1.83	0.61
1:F:61:THR:HA	1:F:94:HIS:CE1	2.35	0.61
1:F:29:ARG:NE	1:F:130:SER:HB2	2.14	0.60
1:F:60:ASP:HA	1:F:74:GLU:OE1	2.01	0.60
1:C:45:GLU:OE2	1:C:92:LYS:NZ	2.37	0.58
1:B:29:ARG:HH11	1:B:29:ARG:CG	2.17	0.58
1:A:2:LYS:NZ	1:A:2:LYS:CD	2.63	0.58
1:A:114:ARG:HG2	1:A:114:ARG:HH11	1.68	0.58
1:B:35:GLN:H	1:B:35:GLN:NE2	2.03	0.57
1:F:31:ASP:O	1:F:34:ILE:HG12	2.05	0.57
1:E:41:GLU:CB	1:E:42:SER:HB2	2.35	0.56
1:D:72:ASN:ND2	1:D:74:GLU:H	2.02	0.56
1:F:129:VAL:HG22	1:F:129:VAL:O	2.06	0.56
1:B:72:ASN:C	1:B:72:ASN:HD22	2.11	0.55
1:B:106:ASN:OD1	1:B:108:SER:HB2	2.08	0.54
1:B:129:VAL:O	1:B:129:VAL:CG2	2.56	0.54
1:A:19:PRO:HB3	1:B:7:TYR:CE1	2.43	0.54
1:C:81:LEU:CD2	1:C:85:HIS:HA	2.38	0.54
1:D:98:ASN:HD21	1:E:85:HIS:CE1	2.26	0.54
1:B:84:ASN:O	1:B:85:HIS:HB2	2.08	0.53
1:C:94:HIS:ND1	1:C:97:LYS:NZ	2.57	0.53
1:E:116:HIS:CE1	1:E:119:GLN:HE21	2.27	0.53
1:E:78:LEU:HD12	1:E:90:ILE:HD11	1.89	0.53
1:F:82:GLU:HG3	1:F:88:THR:CG2	2.39	0.53
1:A:119:GLN:CG	2:A:131:GTQ:HAF	2.33	0.53
1:E:2:LYS:NZ	1:F:31:ASP:OD1	2.41	0.53
1:F:96:GLU:CD	1:F:96:GLU:N	2.55	0.53
1:E:96:GLU:H	1:E:96:GLU:CD	2.12	0.52
1:F:86:TYR:CD1	1:F:125:LEU:HD13	2.44	0.52
1:E:29:ARG:HH11	1:E:29:ARG:CB	2.20	0.52
1:C:65:LEU:HD11	1:C:101:VAL:HG13	1.91	0.52
1:A:72:ASN:HD21	1:A:74:GLU:HB2	1.74	0.52
1:E:82:GLU:HG3	1:E:88:THR:CG2	2.40	0.52
1:F:29:ARG:HE	1:F:130:SER:CB	2.20	0.52
1:E:31:ASP:O	1:E:34:ILE:HG12	2.11	0.51
1:B:78:LEU:HB2	1:B:90:ILE:HG23	1.91	0.51
1:E:109:CYS:SG	1:E:110:LYS:O	2.59	0.51
1:F:20:ASP:OD2	1:F:22:THR:OG1	2.24	0.51
1:E:82:GLU:HG3	1:E:88:THR:HG23	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:109:CYS:SG	1:F:110:LYS:O	2.62	0.51
1:F:57:LEU:HD11	1:F:65:LEU:HD22	1.93	0.50
1:A:112:GLY:N	1:A:113:PRO:CD	2.74	0.50
1:F:59:MET:CB	1:F:59:MET:SD	2.96	0.50
1:F:58:ALA:HB2	1:F:75:CYS:SG	2.52	0.50
1:E:5:LEU:HD22	1:E:34:ILE:HD12	1.95	0.49
1:F:10:ASN:OD1	1:F:105:LYS:NZ	2.45	0.49
1:D:98:ASN:HD21	1:E:85:HIS:HE1	1.61	0.49
1:D:116:HIS:HD2	1:D:117:TYR:O	1.95	0.49
1:A:1:LYS:HB2	1:A:37:GLN:HE22	1.78	0.48
1:A:5:LEU:HD22	1:A:34:ILE:HD12	1.93	0.48
1:C:78:LEU:HB2	1:C:90:ILE:HG23	1.94	0.48
1:E:29:ARG:HB3	1:E:29:ARG:NH1	2.23	0.48
1:E:47:TYR:CE1	1:E:76:LEU:HD13	2.48	0.48
1:A:10:ASN:HB2	1:A:121:ALA:HA	1.95	0.48
1:E:64:LEU:HD12	1:E:111:ARG:HD3	1.94	0.48
1:E:116:HIS:CE1	1:E:119:GLN:NE2	2.81	0.48
1:F:97:LYS:O	1:F:99:TRP:CD1	2.67	0.47
1:F:56:TYR:CE1	1:F:71:PRO:HD3	2.48	0.47
1:F:1:LYS:HA	1:F:2:LYS:HB3	1.95	0.47
1:A:114:ARG:HG2	1:A:114:ARG:NH1	2.29	0.47
1:B:90:ILE:CD1	1:B:98:ASN:HA	2.45	0.47
1:D:72:ASN:HD22	1:D:72:ASN:C	2.19	0.47
1:D:127:LEU:HG	1:F:20:ASP:HB3	1.96	0.47
1:F:82:GLU:HG3	1:F:88:THR:HG23	1.96	0.47
1:A:72:ASN:HD22	1:A:74:GLU:N	2.13	0.47
1:B:86:TYR:CE1	1:B:125:LEU:HB2	2.50	0.47
1:C:90:ILE:HD12	1:C:98:ASN:HA	1.96	0.47
1:D:110:LYS:HD3	1:D:115:THR:HG22	1.97	0.47
1:F:72:ASN:HD22	1:F:74:GLU:N	2.11	0.46
1:A:56:TYR:CE1	1:A:71:PRO:HD3	2.50	0.46
1:D:96:GLU:OE1	1:D:96:GLU:N	2.38	0.46
1:C:14:PHE:CE2	1:C:29:ARG:HG3	2.51	0.46
1:A:8:CYS:HB2	1:A:124:PHE:CE2	2.51	0.46
1:F:129:VAL:O	1:F:129:VAL:CG2	2.64	0.45
1:D:129:VAL:HG13	1:D:129:VAL:O	2.15	0.45
1:E:79:GLU:HG3	1:E:89:TYR:CE2	2.51	0.45
1:F:60:ASP:C	1:F:62:ASP:H	2.20	0.45
1:D:8:CYS:HB2	1:D:124:PHE:CE2	2.51	0.45
1:B:3:PRO:HB3	1:B:51:THR:HG21	1.99	0.45
1:D:29:ARG:HD3	1:F:20:ASP:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:81:LEU:CD2	1:D:85:HIS:HA	2.47	0.44
1:F:100:PHE:HB2	1:F:115:THR:HB	1.99	0.44
1:C:90:ILE:HD11	1:C:95:ALA:HA	1.98	0.44
1:A:66:TYR:OH	1:A:74:GLU:OE1	2.26	0.44
1:C:90:ILE:CD1	1:C:98:ASN:HA	2.47	0.44
1:B:14:PHE:CE2	1:B:29:ARG:HD2	2.53	0.44
1:B:8:CYS:HB2	1:B:124:PHE:CE2	2.53	0.43
1:B:29:ARG:HB2	1:B:29:ARG:NH1	2.25	0.43
1:E:10:ASN:HB2	1:E:121:ALA:HA	2.00	0.43
1:A:89:TYR:HB2	1:A:101:VAL:CG2	2.49	0.43
1:C:104:LYS:HG2	1:C:110:LYS:HG3	2.01	0.43
1:A:105:LYS:HZ1	2:A:131:GTQ:CAH	2.31	0.43
1:F:1:LYS:HA	1:F:2:LYS:O	2.19	0.43
1:E:41:GLU:HG3	1:E:42:SER:HB2	2.01	0.43
1:E:72:ASN:HD22	1:E:74:GLU:H	1.67	0.43
1:D:115:THR:HA	1:D:119:GLN:OE1	2.19	0.43
1:A:72:ASN:HD22	1:A:74:GLU:H	1.66	0.42
1:F:29:ARG:HH21	1:F:130:SER:CB	2.33	0.42
1:F:125:LEU:HD12	1:F:125:LEU:HA	1.86	0.42
1:F:1:LYS:HA	1:F:2:LYS:CB	2.50	0.42
1:F:4:LYS:HD2	1:F:38:LEU:HD12	2.01	0.42
1:E:129:VAL:O	1:E:130:SER:O	2.36	0.42
1:C:2:LYS:HA	1:C:3:PRO:HD2	1.94	0.42
1:F:6:LEU:HD23	1:F:6:LEU:HA	1.90	0.42
1:B:102:GLY:HA2	1:B:121:ALA:O	2.19	0.42
1:E:129:VAL:O	1:E:130:SER:C	2.58	0.42
1:B:90:ILE:HD12	1:B:98:ASN:HA	2.02	0.41
1:D:29:ARG:NH2	1:F:67:GLY:O	2.47	0.41
1:D:47:TYR:CZ	1:D:76:LEU:HD13	2.55	0.41
1:A:72:ASN:ND2	1:A:74:GLU:HB2	2.35	0.41
1:A:90:ILE:HD13	1:A:90:ILE:HG21	1.88	0.41
1:F:72:ASN:ND2	1:F:72:ASN:C	2.72	0.41
1:F:2:LYS:HA	1:F:3:PRO:HD3	1.76	0.41
1:E:14:PHE:O	1:E:25:GLY:HA2	2.21	0.40
1:C:4:LYS:HD2	1:C:38:LEU:HD13	2.04	0.40
1:E:35:GLN:O	1:E:51:THR:OG1	2.39	0.40
1:B:84:ASN:O	1:B:85:HIS:CB	2.69	0.40
1:D:49:LYS:HD3	1:D:56:TYR:CZ	2.56	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	128/130 (98%)	119 (93%)	8 (6%)	1 (1%)	19	17
1	B	128/130 (98%)	122 (95%)	5 (4%)	1 (1%)	19	17
1	C	128/130 (98%)	123 (96%)	3 (2%)	2 (2%)	9	5
1	D	128/130 (98%)	121 (94%)	5 (4%)	2 (2%)	9	5
1	E	128/130 (98%)	121 (94%)	6 (5%)	1 (1%)	19	17
1	F	128/130 (98%)	120 (94%)	5 (4%)	3 (2%)	6	3
All	All	768/780 (98%)	726 (94%)	32 (4%)	10 (1%)	12	8

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	42	SER
1	F	41	GLU
1	B	83	GLU
1	D	43	VAL
1	C	112	GLY
1	F	98	ASN
1	C	41	GLU
1	D	83	GLU
1	F	2	LYS
1	A	43	VAL

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	114/114 (100%)	103 (90%)	11 (10%)	8 6
1	B	114/114 (100%)	101 (89%)	13 (11%)	5 3
1	C	114/114 (100%)	102 (90%)	12 (10%)	7 5
1	D	114/114 (100%)	103 (90%)	11 (10%)	8 6
1	E	114/114 (100%)	98 (86%)	16 (14%)	3 2
1	F	114/114 (100%)	98 (86%)	16 (14%)	3 2
All	All	684/684 (100%)	605 (88%)	79 (12%)	5 3

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LYS
1	A	2	LYS
1	A	32	GLN
1	A	41	GLU
1	A	72	ASN
1	A	76	LEU
1	A	83	GLU
1	A	84	ASN
1	A	90	ILE
1	A	92	LYS
1	A	127	LEU
1	B	1	LYS
1	B	29	ARG
1	B	35	GLN
1	B	64	LEU
1	B	72	ASN
1	B	76	LEU
1	B	81	LEU
1	B	83	GLU
1	B	91	SER
1	B	108	SER
1	B	114	ARG
1	B	127	LEU
1	B	129	VAL
1	C	1	LYS
1	C	2	LYS
1	C	35	GLN
1	C	43	VAL
1	C	72	ASN
1	C	76	LEU

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Mol	Chain	Res	Type
1	C	81	LEU
1	C	83	GLU
1	C	84	ASN
1	C	90	ILE
1	C	116	HIS
1	C	127	LEU
1	D	1	LYS
1	D	32	GLN
1	D	35	GLN
1	D	49	LYS
1	D	64	LEU
1	D	72	ASN
1	D	81	LEU
1	D	98	ASN
1	D	120	LYS
1	D	127	LEU
1	D	130	SER
1	E	1	LYS
1	E	9	SER
1	E	41	GLU
1	E	49	LYS
1	E	51	THR
1	E	64	LEU
1	E	72	ASN
1	E	76	LEU
1	E	80	ARG
1	E	81	LEU
1	E	83	GLU
1	E	96	GLU
1	E	98	ASN
1	E	109	CYS
1	E	120	LYS
1	E	129	VAL
1	F	1	LYS
1	F	37	GLN
1	F	42	SER
1	F	43	VAL
1	F	45	GLU
1	F	64	LEU
1	F	72	ASN
1	F	76	LEU
1	F	80	ARG

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Mol	Chain	Res	Type
1	F	83	GLU
1	F	84	ASN
1	F	96	GLU
1	F	108	SER
1	F	109	CYS
1	F	125	LEU
1	F	129	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	37	GLN
1	A	72	ASN
1	A	84	ASN
1	A	94	HIS
1	B	35	GLN
1	B	69	GLN
1	B	72	ASN
1	B	98	ASN
1	C	35	GLN
1	C	69	GLN
1	C	72	ASN
1	D	35	GLN
1	D	69	GLN
1	D	72	ASN
1	D	85	HIS
1	D	98	ASN
1	D	116	HIS
1	E	13	HIS
1	E	32	GLN
1	E	72	ASN
1	E	119	GLN
1	F	72	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)

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	$\#\|Z\| > 2$
2	GTQ	B	131	-	11,11,11	1.69	1 (9%)	15,15,15	1.64	1 (6%)
2	GTQ	A	131	-	11,11,11	1.33	1 (9%)	15,15,15	1.17	2 (13%)

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
2	GTQ	B	131	-	-	0/4/4/4	0/1/1/1
2	GTQ	A	131	-	-	4/4/4/4	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	131	GTQ	CAK-CAH	4.48	1.59	1.49
2	A	131	GTQ	CAK-CAH	3.81	1.57	1.49

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	131	GTQ	CAJ-CAK-CAH	4.30	124.61	120.03
2	A	131	GTQ	CAJ-CAK-CAH	2.40	122.59	120.03
2	A	131	GTQ	OAB-CAH-CAK	-2.16	116.70	121.94

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	131	GTQ	OAA-CAH-CAK-CAJ
2	A	131	GTQ	OAB-CAH-CAK-CAJ
2	A	131	GTQ	OAA-CAH-CAK-CAG
2	A	131	GTQ	OAB-CAH-CAK-CAG

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	131	GTQ	1	0
2	A	131	GTQ	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	130/130 (100%)	0.81	12 (9%)	9	9	31, 38, 48, 54	0
1	B	130/130 (100%)	0.50	7 (5%)	25	28	28, 37, 46, 100	0
1	C	130/130 (100%)	0.43	7 (5%)	25	28	29, 37, 47, 54	0
1	D	130/130 (100%)	0.41	7 (5%)	25	28	28, 37, 48, 55	0
1	E	130/130 (100%)	0.50	14 (10%)	5	5	29, 37, 47, 57	0
1	F	130/130 (100%)	0.90	14 (10%)	5	5	30, 39, 48, 52	0
All	All	780/780 (100%)	0.59	61 (7%)	13	14	28, 38, 48, 100	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	LYS	7.5
1	D	130	SER	6.8
1	B	130	SER	6.8
1	F	41	GLU	6.7
1	E	41	GLU	6.2
1	D	1	LYS	6.0
1	A	130	SER	5.7
1	F	1	LYS	5.5
1	C	130	SER	5.5
1	C	1	LYS	5.3
1	F	42	SER	5.2
1	F	60	ASP	5.1
1	A	42	SER	4.9
1	A	2	LYS	4.9
1	E	1	LYS	4.4
1	A	61	THR	4.4
1	C	2	LYS	4.3
1	D	129	VAL	4.1
1	B	43	VAL	4.0

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Mol	Chain	Res	Type	RSRZ
1	C	41	GLU	3.8
1	F	64	LEU	3.6
1	F	105	LYS	3.6
1	B	129	VAL	3.4
1	F	43	VAL	3.4
1	B	1	LYS	3.4
1	F	73	GLU	3.4
1	E	9	SER	3.2
1	C	43	VAL	3.1
1	C	129	VAL	3.0
1	B	2	LYS	3.0
1	F	56	TYR	3.0
1	D	2	LYS	2.9
1	A	43	VAL	2.9
1	F	130	SER	2.9
1	A	40	ALA	2.9
1	E	107	GLY	2.8
1	E	116	HIS	2.8
1	B	83	GLU	2.8
1	F	118	GLY	2.7
1	E	40	ALA	2.6
1	E	73	GLU	2.6
1	D	41	GLU	2.5
1	E	42	SER	2.5
1	D	3	PRO	2.5
1	E	29	ARG	2.4
1	C	105	LYS	2.4
1	B	61	THR	2.4
1	F	99	TRP	2.4
1	F	106	ASN	2.4
1	E	106	ASN	2.3
1	A	63	GLY	2.3
1	A	41	GLU	2.2
1	A	72	ASN	2.2
1	D	84	ASN	2.2
1	E	64	LEU	2.2
1	E	10	ASN	2.2
1	F	81	LEU	2.1
1	A	3	PRO	2.1
1	E	130	SER	2.1
1	E	11	GLY	2.1
1	A	129	VAL	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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	GTQ	B	131	11/11	0.78	0.24	60,62,63,64	3
2	GTQ	A	131	11/11	0.87	0.18	55,57,59,59	3

6.5 Other polymers (i)

There are no such residues in this entry.