Full wwPDB X-ray Structure Validation Report (i)

Jul 26, 2023 – 01:56 AM EDT

PDB ID	:	1AVF
Title	:	ACTIVATION INTERMEDIATE 2 OF HUMAN GASTRICSIN FROM HUMAN STOMACH
Authors	:	Khan, A.R.; Cherney, M.M.; Tarasova, N.I.; James, M.N.G.
Deposited on	:	1997-09-16
Resolution	:	2.36 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 (1)) were used in the production of this report:

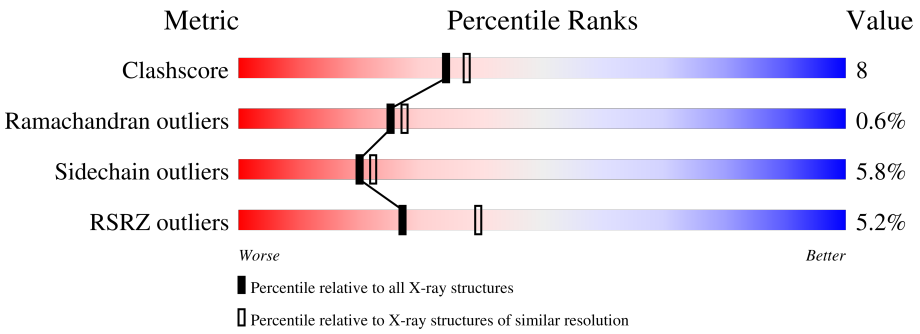
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.34
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.34

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 2.36 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))
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	P	26	8% 69% 12% 19%
1	Q	26	4% 73% 8% • 15%
2	A	329	4% 74% 21% • •
2	J	329	6% 76% 20% • • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5686 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 GASTRICSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	P	21	168	110	30	27	1	0	0	0
1	Q	22	176	116	31	28	1	0	0	0

- Molecule 2 is a protein called GASTRICSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	A	322	2431	1553	377	489	12	0	0	2
2	J	323	2443	1562	377	492	12	0	0	1

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
3	A	1	1	1	0	0

- Molecule 4 is water.

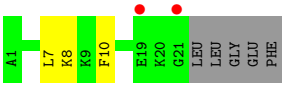
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	P	13	13	13	0	0
4	A	226	226	226	0	0
4	Q	12	12	12	0	0
4	J	216	216	216	0	0

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.

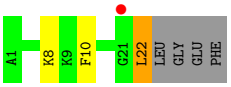
- Molecule 1: GASTRICSIN

Chain P:



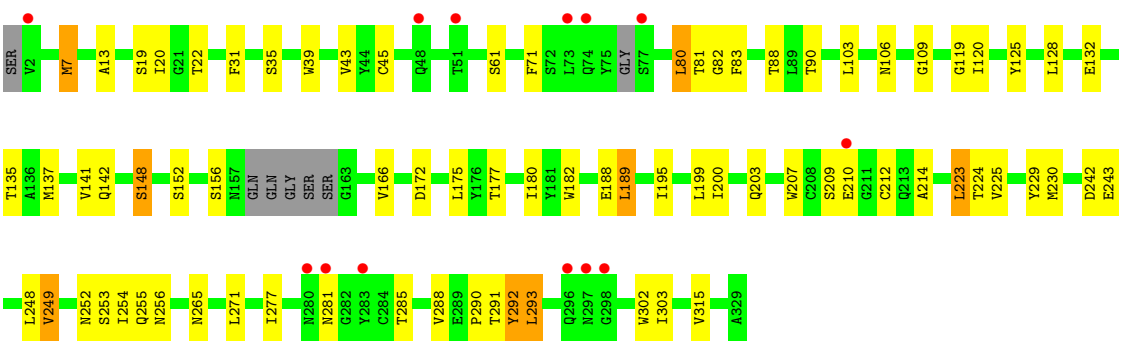
- Molecule 1: GASTRICSIN

Chain Q:



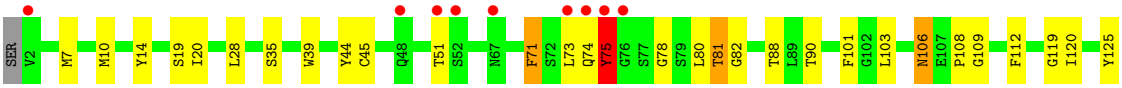
- Molecule 2: GASTRICSIN

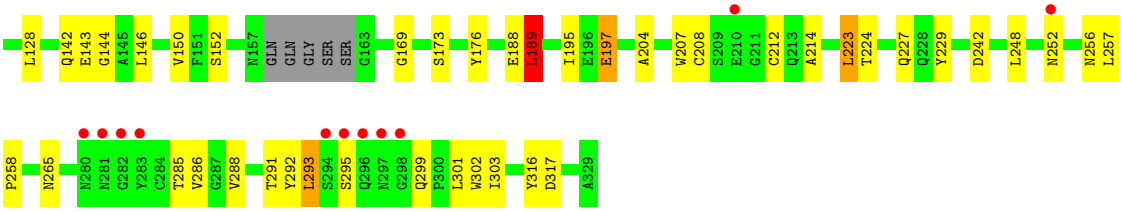
Chain A:



- Molecule 2: GASTRICSIN

Chain J:




L128
Q142
E143
G144
A145
L146
V150
F151
S152
N157
GLN
GLN
GLY
SER
SER
G163
G169
S173
Y176
E188
L189
I195
E196
E197
A204
W207
C208
S209
E210
G211
C212
Q213
A214
L223
T224
Q227
Q228
Y229
D242
L248
N252
N256
L257
P258
N265
N280
N281
G282
Y283
C284
T285
V286
G287
V288
T291
Y292
L293
S294
S295
Q296
N297
G298
Q299
P300
L301
W302
I303
Y316
D317
A329

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	156.26Å 50.41Å 125.18Å 90.00° 117.62° 90.00°	Depositor
Resolution (Å)	20.00 – 2.36 19.77 – 2.36	Depositor EDS
% Data completeness (in resolution range)	95.3 (20.00-2.36) 95.7 (19.77-2.36)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.12	Depositor
$< I/\sigma(I) >$ [1]	2.45 (at 2.35Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.227 , 0.282 0.188 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	26.5	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 66.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5686	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: NA

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	P	0.56	0/169	0.77	0/221
1	Q	0.60	0/177	0.96	1/232 (0.4%)
2	A	0.56	0/2494	0.78	1/3409 (0.0%)
2	J	0.53	0/2508	0.76	1/3429 (0.0%)
All	All	0.55	0/5348	0.78	3/7291 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	Q	22	LEU	N-CA-C	5.58	126.06	111.00
2	J	189	LEU	N-CA-C	-5.05	97.36	111.00
2	A	189	LEU	N-CA-C	-5.02	97.44	111.00

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	P	168	0	199	4	0
1	Q	176	0	210	1	0
2	A	2431	0	2255	40	17
2	J	2443	0	2267	42	15

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
4	A	226	0	0	7	31
4	J	216	0	0	3	11
4	P	13	0	0	1	0
4	Q	12	0	0	0	1
All	All	5686	0	4931	85	42

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 (85) 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)
2:A:203:GLN:HG3	4:A:527:HOH:O	1.66	0.94
2:J:248:LEU:HD23	2:J:285:THR:HG22	1.67	0.76
2:A:7:MET:HG3	2:A:109:GLY:HA3	1.69	0.75
2:J:7:MET:HG3	2:J:109:GLY:HA3	1.72	0.71
2:J:223:LEU:HB3	2:J:288:VAL:HG12	1.73	0.69
2:A:148:SER:HB3	4:A:485:HOH:O	1.93	0.68
2:A:180:ILE:HD13	2:A:315:VAL:HG21	1.77	0.66
2:A:200:ILE:O	2:A:203:GLN:HG2	1.95	0.65
2:A:156:SER:HB2	4:A:526:HOH:O	1.98	0.63
2:A:224:THR:HG22	2:A:291:THR:HB	1.83	0.60
2:J:82:GLY:HA3	2:J:103:LEU:O	2.02	0.59
2:J:214:ALA:HA	2:J:302:TRP:O	2.01	0.59
2:J:35:SER:HB3	4:J:334:HOH:O	2.03	0.57
2:J:224:THR:HG22	2:J:291:THR:HB	1.87	0.57
2:A:135:THR:HG22	2:A:137:MET:H	1.70	0.56
2:A:224:THR:CG2	2:A:291:THR:HB	2.37	0.55
2:J:224:THR:HB	2:J:303:ILE:HB	1.88	0.55
2:A:182:TRP:HB2	4:A:524:HOH:O	2.05	0.55
1:P:7:LEU:HD21	2:A:166:VAL:HG23	1.89	0.55
2:J:119:GLY:C	2:J:120:ILE:HD12	2.28	0.53
2:J:101:PHE:CE2	2:J:103:LEU:HD13	2.43	0.52
2:A:195:ILE:HB	2:A:212:CYS:HB2	1.92	0.52
2:A:249:VAL:HG21	2:A:277:ILE:HD12	1.92	0.52
2:J:101:PHE:HE2	2:J:103:LEU:HD13	1.75	0.52
2:A:7:MET:HG3	2:A:109:GLY:CA	2.39	0.51
2:J:197:GLU:HA	2:J:208:CYS:HB2	1.92	0.51
2:J:224:THR:CG2	2:J:291:THR:HB	2.40	0.51
1:P:8:LYS:HD2	4:P:39:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:210:GLU:HB3	4:A:492:HOH:O	2.10	0.50
2:J:173:SER:HA	2:J:176:TYR:CE1	2.47	0.50
2:J:75:TYR:C	2:J:75:TYR:CD1	2.85	0.50
2:A:225:VAL:HG22	2:A:302:TRP:CZ3	2.46	0.50
1:P:8:LYS:HB2	1:P:10:PHE:CE1	2.47	0.49
2:A:135:THR:HG22	2:A:137:MET:N	2.27	0.49
2:J:125:TYR:CE2	2:J:188:GLU:HB3	2.47	0.49
2:J:108:PRO:HD2	2:J:112:PHE:CD2	2.47	0.48
2:J:71:PHE:O	2:J:81:THR:HA	2.13	0.48
2:J:195:ILE:HD11	2:J:214:ALA:HB2	1.96	0.48
2:A:31:PHE:CD2	2:A:166:VAL:HG21	2.50	0.47
2:A:22:THR:O	2:A:61:SER:HA	2.15	0.47
2:A:265:ASN:HB2	4:A:398:HOH:O	2.15	0.47
2:J:224:THR:HG21	2:J:303:ILE:HD12	1.97	0.47
2:J:150:VAL:HG22	2:J:317:ASP:HA	1.97	0.47
2:J:19:SER:HB2	2:J:90:THR:HB	1.98	0.46
1:P:10:PHE:O	2:A:13:ALA:HB1	2.16	0.46
2:A:224:THR:HB	2:A:303:ILE:HB	1.97	0.46
2:A:248:LEU:HD23	2:A:285:THR:HG22	1.97	0.46
2:A:230:MET:HG3	2:A:290:PRO:HB3	1.99	0.45
2:A:223:LEU:HB3	2:A:288:VAL:HG12	1.99	0.45
2:J:39:TRP:CZ2	2:J:120:ILE:HD13	2.53	0.44
2:J:195:ILE:HB	2:J:212:CYS:HB2	1.98	0.44
2:A:82:GLY:HA3	2:A:103:LEU:O	2.17	0.44
2:J:74:GLN:O	2:J:78:GLY:O	2.35	0.44
2:J:316:TYR:CD1	2:J:316:TYR:N	2.86	0.44
2:A:195:ILE:HD11	2:A:214:ALA:HB2	2.00	0.44
2:A:292:TYR:CD1	2:A:292:TYR:N	2.86	0.44
2:A:207:TRP:CH2	2:A:229:TYR:HB3	2.53	0.43
2:J:45:CYS:SG	2:J:106:ASN:HA	2.58	0.43
2:J:197:GLU:OE2	2:J:204:ALA:HB1	2.19	0.43
2:A:39:TRP:CE2	2:A:80:LEU:HD21	2.54	0.43
2:A:19:SER:HB2	2:A:90:THR:HB	2.01	0.43
2:A:20:ILE:HA	2:A:88:THR:O	2.19	0.42
2:J:125:TYR:HB2	2:J:189:LEU:O	2.19	0.42
2:A:125:TYR:CE2	2:A:188:GLU:HB3	2.55	0.42
2:J:207:TRP:HB3	2:J:229:TYR:CE2	2.54	0.42
2:J:20:ILE:HA	2:J:88:THR:O	2.20	0.42
2:A:80:LEU:HD12	2:A:80:LEU:HA	1.93	0.42
2:J:223:LEU:O	2:J:288:VAL:HA	2.20	0.42
2:J:293:LEU:HD23	2:J:301:LEU:HD23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:189:LEU:HD23	2:J:189:LEU:HA	1.95	0.41
2:J:146:LEU:HB3	2:J:169:GLY:HA2	2.01	0.41
2:A:45:CYS:SG	2:A:106:ASN:HA	2.60	0.41
2:A:292:TYR:N	2:A:292:TYR:HD1	2.18	0.41
2:J:292:TYR:N	2:J:292:TYR:CD1	2.89	0.41
2:A:172:ASP:HB3	2:A:175:LEU:HD12	2.03	0.40
1:Q:8:LYS:HB3	1:Q:10:PHE:CE1	2.55	0.40
2:J:14:TYR:HD2	4:J:365:HOH:O	2.04	0.40
2:J:265:ASN:HB2	4:J:399:HOH:O	2.21	0.40
2:A:293:LEU:HD12	2:A:293:LEU:HA	1.84	0.40
2:J:73:LEU:HD12	2:J:80:LEU:HD23	2.04	0.40
2:J:295:SER:HA	2:J:299:GLN:O	2.21	0.40
2:A:119:GLY:C	2:A:120:ILE:HD12	2.42	0.40
2:J:28:LEU:HD23	2:J:28:LEU:HA	1.76	0.40
2:J:286:VAL:HG12	2:J:288:VAL:HG22	2.04	0.40
2:A:35:SER:HB3	4:A:338:HOH:O	2.20	0.40

All (42) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:475:HOH:O	4:A:485:HOH:O[2_556]	0.11	2.09
4:A:341:HOH:O	4:A:457:HOH:O[2_556]	0.46	1.74
4:J:475:HOH:O	4:J:519:HOH:O[4_445]	0.47	1.73
2:J:257:LEU:C	4:J:530:HOH:O[4_455]	0.72	1.48
4:A:444:HOH:O	4:A:538:HOH:O[2_556]	0.75	1.45
2:A:242:ASP:O	4:A:539:HOH:O[4_556]	0.76	1.44
2:J:257:LEU:CA	4:J:530:HOH:O[4_455]	0.80	1.40
2:A:142:GLN:N	4:A:529:HOH:O[2_556]	0.87	1.33
2:A:242:ASP:C	4:A:539:HOH:O[4_556]	0.93	1.27
2:A:177:THR:CB	4:A:545:HOH:O[4_546]	1.35	0.85
2:A:255:GLN:CA	4:A:544:HOH:O[4_546]	1.36	0.84
2:A:142:GLN:CA	4:A:529:HOH:O[2_556]	1.38	0.82
2:J:44:TYR:CD1	4:A:516:HOH:O[1_545]	1.42	0.78
2:J:242:ASP:O	4:J:524:HOH:O[4_445]	1.42	0.78
2:A:243:GLU:N	4:A:539:HOH:O[4_556]	1.51	0.69
2:A:141:VAL:C	4:A:529:HOH:O[2_556]	1.53	0.67
2:J:144:GLY:N	4:A:524:HOH:O[2_556]	1.66	0.54
2:J:257:LEU:O	4:J:530:HOH:O[4_455]	1.67	0.53
2:A:256:ASN:N	4:A:530:HOH:O[4_546]	1.69	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:142:GLN:O	4:A:524:HOH:O[2_556]	1.82	0.38
4:A:550:HOH:O	4:Q:89:HOH:O[2_556]	1.82	0.38
2:J:258:PRO:N	4:J:530:HOH:O[4_455]	1.82	0.38
2:J:257:LEU:N	4:J:530:HOH:O[4_455]	1.84	0.36
4:A:459:HOH:O	4:A:555:HOH:O[4_546]	1.88	0.32
2:J:256:ASN:O	4:J:502:HOH:O[4_455]	1.89	0.31
2:A:255:GLN:N	4:A:530:HOH:O[4_546]	1.90	0.30
2:J:142:GLN:C	4:A:524:HOH:O[2_556]	1.92	0.28
2:J:257:LEU:CB	4:J:530:HOH:O[4_455]	1.93	0.27
2:A:243:GLU:CA	4:A:539:HOH:O[4_556]	1.97	0.23
4:A:350:HOH:O	4:A:509:HOH:O[4_546]	1.98	0.22
2:J:143:GLU:C	4:A:524:HOH:O[2_556]	2.01	0.19
2:A:255:GLN:N	4:A:544:HOH:O[4_546]	2.03	0.17
2:A:253:SER:O	4:A:530:HOH:O[4_546]	2.08	0.12
4:A:451:HOH:O	4:A:530:HOH:O[4_546]	2.08	0.12
2:A:177:THR:CG2	4:A:545:HOH:O[4_546]	2.09	0.11
2:J:242:ASP:N	4:J:524:HOH:O[4_445]	2.09	0.11
2:A:255:GLN:CB	4:A:510:HOH:O[4_546]	2.10	0.10
4:J:392:HOH:O	4:J:525:HOH:O[4_445]	2.13	0.07
2:J:44:TYR:CE1	4:A:516:HOH:O[1_545]	2.14	0.06
4:A:350:HOH:O	4:A:555:HOH:O[4_546]	2.15	0.05
2:A:177:THR:CA	4:A:545:HOH:O[4_546]	2.16	0.04
2:A:141:VAL:O	4:A:529:HOH:O[2_556]	2.17	0.03

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	P	19/26 (73%)	19 (100%)	0	0	100	100
1	Q	20/26 (77%)	20 (100%)	0	0	100	100
2	A	316/329 (96%)	307 (97%)	7 (2%)	2 (1%)	25	27
2	J	319/329 (97%)	304 (95%)	13 (4%)	2 (1%)	25	27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	674/710 (95%)	650 (96%)	20 (3%)	4 (1%)	25	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	J	75	TYR
2	A	281	ASN
2	A	254	ILE
2	J	51	THR

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	P	19/23 (83%)	19 (100%)	0	100	100
1	Q	20/23 (87%)	19 (95%)	1 (5%)	24	28
2	A	263/274 (96%)	244 (93%)	19 (7%)	14	14
2	J	264/274 (96%)	251 (95%)	13 (5%)	25	29
All	All	566/594 (95%)	533 (94%)	33 (6%)	20	22

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	7	MET
2	A	43	VAL
2	A	71	PHE
2	A	80	LEU
2	A	81	THR
2	A	83	PHE
2	A	128	LEU
2	A	132	GLU
2	A	148	SER
2	A	152	SER
2	A	189	LEU

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Mol	Chain	Res	Type
2	A	199	LEU
2	A	209	SER
2	A	223	LEU
2	A	249	VAL
2	A	252	ASN
2	A	271	LEU
2	A	292	TYR
2	A	293	LEU
1	Q	22	LEU
2	J	10	MET
2	J	71	PHE
2	J	75	TYR
2	J	81	THR
2	J	106	ASN
2	J	128	LEU
2	J	152	SER
2	J	189	LEU
2	J	197	GLU
2	J	223	LEU
2	J	227	GLN
2	J	252	ASN
2	J	293	LEU

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

Mol	Chain	Res	Type
2	A	74	GLN
2	A	95	GLN
2	J	46	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is 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.

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	P	21/26 (80%)	-0.03	2 (9%)	8	13	4, 19, 51, 53	0
1	Q	22/26 (84%)	-0.08	1 (4%)	33	46	7, 22, 49, 51	0
2	A	322/329 (97%)	-0.25	13 (4%)	38	51	2, 12, 39, 64	0
2	J	323/329 (98%)	-0.15	20 (6%)	20	29	2, 14, 45, 71	0
All	All	688/710 (96%)	-0.19	36 (5%)	27	39	2, 13, 46, 71	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	75	TYR	6.9
2	J	74	GLN	5.8
2	A	281	ASN	5.7
2	J	282	GLY	5.0
1	Q	21	GLY	4.6
2	A	74	GLN	4.6
2	J	297	ASN	4.5
2	J	73	LEU	4.2
2	J	76	GLY	4.2
2	J	280	ASN	4.1
2	J	48	GLN	4.0
2	A	297	ASN	3.6
2	J	2	VAL	3.5
2	A	283	TYR	3.4
2	J	281	ASN	3.3
2	A	280	ASN	3.3
2	A	51	THR	3.2
2	J	298	GLY	3.1
2	A	298	GLY	3.1
1	P	19	GLU	2.9
2	A	48	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
1	P	21	GLY	2.7
2	A	210	GLU	2.6
2	J	295	SER	2.4
2	J	283	TYR	2.3
2	J	210	GLU	2.3
2	J	67	ASN	2.3
2	J	252	ASN	2.2
2	J	296	GLN	2.2
2	A	2	VAL	2.2
2	A	73	LEU	2.1
2	A	296	GLN	2.1
2	J	51	THR	2.1
2	A	77	SER	2.1
2	J	52	SER	2.0
2	J	294	SER	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
3	NA	A	330	1/1	0.83	0.07	7,7,7,7	0

6.5 Other polymers (i)

There are no such residues in this entry.