



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:52 PM GMT

PDB ID : 1CZI  
Title : CHYMOSIN COMPLEX WITH THE INHIBITOR CP-113972  
Authors : Groves, M.R.; Dhanaraj, V.; Pitts, J.E.; Badasso, M.; Hoover, D.; Nugent, P.;  
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Deposited on : 1997-01-15  
Resolution : 2.30 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

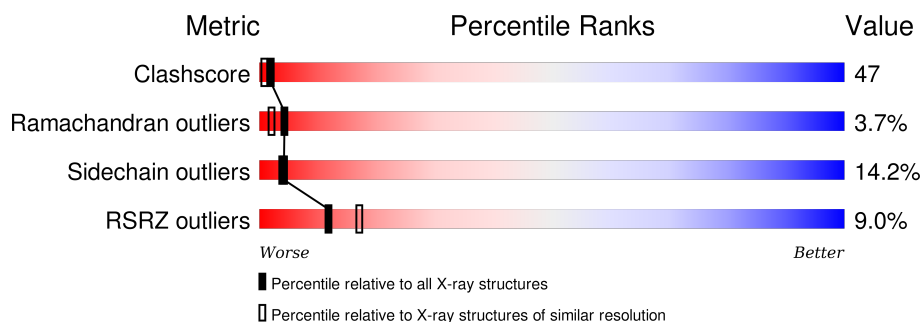
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	E	323	
2	P	4	

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
2	PHI	P	3	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2745 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 CHYMOSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	323	Total	C	N	O	S	0	0	0
			2511	1598	401	498	14			

- Molecule 2 is a protein called CP-113972 (NORSTATINE-S-METHYL CYSTEINE-iodo-PHENYLALANINE-PROLINE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	4	Total	C	I	N	O	S	0	0
			43	31	1	4	6	1		

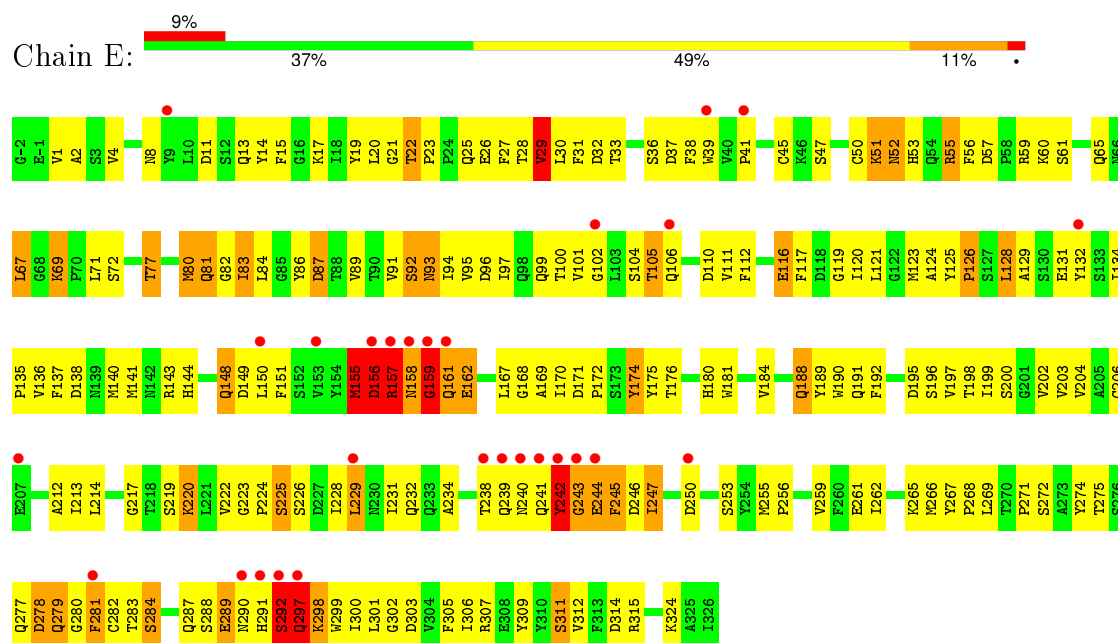
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	191	Total	O	0	0
			191	191		

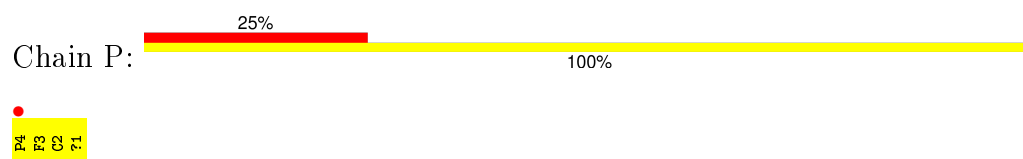
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: CHYMOSIN



#### • Molecule 2: CP-113972 (NORSTATINE-S-METHYL CYSTEINE-IODO-PHENYLALANINE-PROLINE)



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	132.78 Å   132.78 Å   81.95 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	9.97 – 2.30 21.01 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.0 (9.97-2.30) 97.6 (21.01-2.30)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.50 (at 2.30 Å)	Xtriage
Refinement program	RESTRAIN	Depositor
R, $R_{free}$	0.195 , (Not available) 0.266 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	14.5	Xtriage
Anisotropy	0.227	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 19.7	EDS
Estimated twinning fraction	0.023 for -1/3*h+1/3*k+4/3*l,-k,2/3*h+1/3*k+1/3*l 0.035 for -2/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+4/3*l,-1/3*h+1/3*k+1/3*l 0.017 for -h,1/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+1/3*l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 12125 reflections	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	2745	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	0.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: PHI, SMC, NOR

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	E	0.46	0/2574	1.40	18/3506 (0.5%)
2	P	0.24	0/7	0.89	0/8
All	All	0.46	0/2581	1.40	18/3514 (0.5%)

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	4

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	159	GLY	C-N-CA	16.02	161.74	121.70
1	E	156	ASP	C-N-CA	10.95	149.09	121.70
1	E	278	ASP	CB-CG-OD1	-7.57	111.49	118.30
1	E	309	TYR	CB-CG-CD1	-7.48	116.51	121.00
1	E	157	ARG	C-N-CA	7.44	140.30	121.70
1	E	55	ARG	NE-CZ-NH1	-7.13	116.73	120.30
1	E	174	TYR	CB-CG-CD1	-6.53	117.08	121.00
1	E	189	TYR	CB-CG-CD1	6.16	124.70	121.00
1	E	309	TYR	CB-CG-CD2	6.14	124.68	121.00
1	E	155	MET	CA-C-N	5.93	130.24	117.20
1	E	189	TYR	CB-CG-CD2	-5.88	117.47	121.00
1	E	188	GLN	O-C-N	5.80	131.98	122.70
1	E	87	ASP	CB-CG-OD2	5.56	123.30	118.30
1	E	29	VAL	CA-CB-CG2	5.55	119.23	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	289	GLU	O-C-N	5.49	131.48	122.70
1	E	242	TYR	C-N-CA	5.36	133.56	122.30
1	E	29	VAL	CG1-CB-CG2	5.21	119.23	110.90
1	E	37	ASP	CB-CG-OD2	5.13	122.92	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	143	ARG	Sidechain
1	E	155	MET	Peptide
1	E	157	ARG	Sidechain
1	E	159	GLY	Peptide

## 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	E	2511	0	2382	228	6
2	P	43	0	43	20	0
3	E	191	0	0	25	4
All	All	2745	0	2425	233	10

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 47.

All (233) 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:P:3:PHI:I	2:P:3:PHI:CZ	1.36	1.74
2:P:3:PHI:I	2:P:3:PHI:CE2	2.42	1.36
2:P:3:PHI:I	2:P:3:PHI:CE1	2.44	1.35
1:E:77:THR:HG22	1:E:77:THR:O	1.54	1.03
1:E:243:GLY:O	1:E:244:GLU:HB2	1.58	1.02
1:E:297:GLN:HG3	1:E:298:LYS:N	1.69	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:278:ASP:HB3	1:E:281:PHE:HD1	1.25	1.01
1:E:91:VAL:O	1:E:92:SER:HB3	1.60	1.00
1:E:25:GLN:HE22	1:E:57:ASP:H	1.17	0.91
1:E:278:ASP:HB3	1:E:281:PHE:CD1	2.07	0.90
1:E:214:LEU:HD21	1:E:311:SER:HB3	1.55	0.88
1:E:77:THR:CG2	1:E:77:THR:O	2.27	0.82
1:E:171:ASP:HB3	1:E:174:TYR:CD2	2.15	0.81
1:E:176:THR:CG2	1:E:324:LYS:HB3	2.11	0.81
1:E:224:PRO:HB3	1:E:292:SER:OG	1.83	0.79
1:E:13:GLN:NE2	2:P:3:PHI:I	2.85	0.79
1:E:156:ASP:HB2	1:E:157:ARG:HD3	1.63	0.79
1:E:292:SER:O	1:E:297:GLN:HB3	1.83	0.79
1:E:157:ARG:HA	1:E:161:GLN:HE21	1.49	0.78
1:E:277:GLN:NE2	1:E:280:GLY:H	1.84	0.76
1:E:181:TRP:HH2	1:E:312:VAL:HG11	1.49	0.76
1:E:2:ALA:HB2	1:E:94:ILE:HG13	1.69	0.75
1:E:222:VAL:HG12	1:E:289:GLU:HG3	1.69	0.74
1:E:171:ASP:HB3	1:E:174:TYR:HD2	1.50	0.74
1:E:195:ASP:HB2	1:E:261:GLU:HG2	1.70	0.73
1:E:240:ASN:CG	1:E:241:GLN:H	1.92	0.72
1:E:265:LYS:HB3	3:E:613:HOH:O	1.88	0.72
1:E:71:LEU:HD22	1:E:102:GLY:HA3	1.71	0.71
1:E:67:LEU:HG	3:E:438:HOH:O	1.91	0.71
1:E:220:LYS:CE	2:P:4:PRO:HG3	2.22	0.70
1:E:150:LEU:HB3	3:E:549:HOH:O	1.89	0.70
1:E:222:VAL:CG1	1:E:289:GLU:HG3	2.22	0.70
1:E:220:LYS:CD	2:P:4:PRO:HG3	2.22	0.70
1:E:156:ASP:CB	1:E:157:ARG:HD3	2.22	0.70
1:E:30:LEU:HD11	3:E:594:HOH:O	1.91	0.69
1:E:176:THR:HG22	1:E:324:LYS:HB3	1.74	0.69
1:E:297:GLN:CG	1:E:298:LYS:N	2.52	0.69
1:E:71:LEU:CD2	1:E:102:GLY:HA3	2.22	0.69
1:E:181:TRP:CH2	1:E:312:VAL:HG11	2.28	0.68
1:E:259:VAL:HG11	1:E:266:MET:HE3	1.74	0.68
2:P:3:PHI:I	2:P:3:PHI:HE2	2.61	0.68
1:E:150:LEU:HD23	1:E:314:ASP:HA	1.74	0.68
1:E:220:LYS:HG2	2:P:4:PRO:HB3	1.75	0.67
1:E:277:GLN:HE21	1:E:280:GLY:H	1.42	0.67
1:E:128:LEU:CD1	1:E:188:GLN:HB3	2.25	0.67
1:E:128:LEU:HD13	1:E:188:GLN:C	2.17	0.66
1:E:4:VAL:HG23	3:E:557:HOH:O	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:161:GLN:C	1:E:162:GLU:HG2	2.16	0.65
1:E:19:TYR:CE1	1:E:26:GLU:HB2	2.32	0.65
1:E:111:VAL:HB	3:E:532:HOH:O	1.97	0.65
1:E:176:THR:HG23	1:E:324:LYS:HB3	1.80	0.64
1:E:128:LEU:HG	3:E:581:HOH:O	1.96	0.64
1:E:269:LEU:HD13	3:E:491:HOH:O	1.96	0.64
1:E:91:VAL:O	1:E:92:SER:CB	2.34	0.63
1:E:156:ASP:HB3	1:E:307:ARG:HG2	1.80	0.63
1:E:259:VAL:HA	1:E:268:PRO:HA	1.80	0.62
1:E:229:LEU:HD22	3:E:597:HOH:O	1.99	0.62
1:E:2:ALA:HB2	1:E:94:ILE:CD1	2.29	0.62
1:E:51:LYS:HA	1:E:55:ARG:NH1	2.13	0.62
1:E:157:ARG:C	1:E:161:GLN:HB2	2.20	0.62
1:E:22:THR:HA	1:E:23:PRO:C	2.21	0.61
1:E:228:ILE:O	1:E:232:GLN:HG2	2.00	0.61
1:E:156:ASP:CG	1:E:157:ARG:HD3	2.21	0.61
1:E:21:GLY:HA2	1:E:87:ASP:OD1	2.00	0.61
1:E:303:ASP:O	1:E:307:ARG:HB2	2.00	0.61
1:E:155:MET:O	1:E:156:ASP:HB3	1.99	0.61
1:E:2:ALA:HB2	1:E:94:ILE:CG1	2.31	0.61
1:E:121:LEU:C	1:E:121:LEU:HD23	2.21	0.60
1:E:206:CYS:HA	3:E:590:HOH:O	2.02	0.60
1:E:57:ASP:HB3	1:E:60:LYS:HD2	1.83	0.60
1:E:157:ARG:HB2	1:E:158:ASN:OD1	2.01	0.60
1:E:126:PRO:HG3	1:E:138:ASP:OD2	2.03	0.59
2:P:3:PHI:I	2:P:3:PHI:HE1	2.63	0.59
1:E:39:TRP:CE2	1:E:80:MET:HG2	2.37	0.59
1:E:289:GLU:HG2	3:E:507:HOH:O	2.02	0.59
1:E:241:GLN:O	1:E:242:TYR:HB2	2.03	0.59
1:E:240:ASN:CG	1:E:241:GLN:N	2.55	0.58
1:E:77:THR:HG21	2:P:3:PHI:HA	1.84	0.58
1:E:95:VAL:HG12	1:E:97:ILE:HD12	1.84	0.58
1:E:214:LEU:CD2	1:E:311:SER:HB3	2.33	0.58
1:E:8:ASN:OD1	1:E:11:ASP:HA	2.03	0.58
1:E:65:GLN:HB3	1:E:86:TYR:HB2	1.85	0.58
1:E:15:PHE:CE1	1:E:116:GLU:HB3	2.40	0.57
1:E:161:GLN:O	1:E:162:GLU:HG2	2.04	0.57
1:E:162:GLU:HB2	3:E:377:HOH:O	2.05	0.56
1:E:222:VAL:HA	1:E:287:GLN:O	2.05	0.56
1:E:197:VAL:O	1:E:203:VAL:HA	2.05	0.55
1:E:225:SER:HA	1:E:288:SER:HB2	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:274:TYR:O	1:E:274:TYR:HD1	1.90	0.55
1:E:155:MET:O	1:E:307:ARG:HG2	2.07	0.54
1:E:13:GLN:OE1	1:E:116:GLU:HB2	2.07	0.54
1:E:150:LEU:HD13	3:E:549:HOH:O	2.06	0.54
1:E:29:VAL:HG11	1:E:121:LEU:HB2	1.89	0.54
1:E:80:MET:C	1:E:80:MET:SD	2.86	0.54
1:E:51:LYS:HA	1:E:55:ARG:HH11	1.71	0.54
1:E:156:ASP:HB2	1:E:157:ARG:CD	2.35	0.54
1:E:155:MET:O	1:E:156:ASP:CB	2.55	0.54
1:E:67:LEU:CD2	1:E:86:TYR:CE2	2.91	0.54
1:E:148:GLN:HG3	1:E:168:GLY:O	2.08	0.54
1:E:297:GLN:HG3	1:E:298:LYS:H	1.68	0.54
1:E:213:ILE:HD13	3:E:593:HOH:O	2.08	0.54
1:E:120:ILE:HD11	2:P:1:NOR:HE22	1.89	0.53
1:E:156:ASP:CG	1:E:157:ARG:HH11	2.11	0.53
1:E:291:HIS:O	1:E:292:SER:HB2	2.07	0.53
1:E:198:THR:HA	1:E:202:VAL:O	2.09	0.53
1:E:190:TRP:O	1:E:213:ILE:HA	2.08	0.53
1:E:324:LYS:HG2	3:E:621:HOH:O	2.08	0.53
1:E:144:HIS:O	1:E:144:HIS:CD2	2.61	0.52
1:E:150:LEU:HD23	1:E:314:ASP:CA	2.40	0.52
1:E:71:LEU:HD23	1:E:80:MET:HE3	1.91	0.52
1:E:31:PHE:CE1	1:E:121:LEU:HD13	2.44	0.52
1:E:199:ILE:HD12	1:E:234:ALA:HB1	1.91	0.52
1:E:271:PRO:HA	1:E:274:TYR:CE2	2.44	0.52
1:E:141:MET:SD	1:E:149:ASP:HB3	2.49	0.52
1:E:184:VAL:HG13	1:E:191:GLN:O	2.10	0.52
1:E:128:LEU:HD13	1:E:188:GLN:O	2.10	0.52
1:E:213:ILE:HG13	1:E:213:ILE:O	2.08	0.52
1:E:126:PRO:HD3	1:E:315:ARG:HH12	1.75	0.52
1:E:83:ILE:HD13	1:E:83:ILE:N	2.25	0.52
1:E:38:PHE:CE2	1:E:101:VAL:HB	2.45	0.51
1:E:277:GLN:HE21	1:E:279:GLN:HA	1.75	0.51
1:E:212:ALA:HA	1:E:299:TRP:O	2.11	0.51
1:E:247:ILE:O	1:E:282:CYS:HB2	2.11	0.51
1:E:196:SER:HB2	1:E:203:VAL:HG13	1.93	0.51
1:E:128:LEU:HD13	1:E:188:GLN:HB3	1.93	0.51
1:E:213:ILE:HG21	3:E:593:HOH:O	2.10	0.51
1:E:269:LEU:HD11	1:E:305:PHE:HD1	1.75	0.50
1:E:125:TYR:HB3	1:E:126:PRO:HD2	1.92	0.50
1:E:223:GLY:O	1:E:288:SER:HA	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:67:LEU:HD22	1:E:86:TYR:CD2	2.46	0.50
1:E:36:SER:HB3	1:E:135:PRO:HG3	1.93	0.50
1:E:45:CYS:HB2	1:E:104:SER:O	2.12	0.50
1:E:100:THR:HG23	3:E:538:HOH:O	2.12	0.49
1:E:300:ILE:HD13	2:P:1:NOR:HM12	1.94	0.49
1:E:19:TYR:HA	1:E:25:GLN:O	2.12	0.49
1:E:277:GLN:HG3	1:E:278:ASP:N	2.27	0.49
1:E:72:SER:HB2	1:E:81:GLN:HE21	1.78	0.49
1:E:297:GLN:HG2	3:E:502:HOH:O	2.13	0.48
1:E:14:TYR:O	1:E:30:LEU:HD12	2.12	0.48
1:E:150:LEU:HD22	1:E:312:VAL:CG1	2.43	0.48
1:E:69:LYS:O	1:E:84:LEU:HB2	2.13	0.48
1:E:86:TYR:CE1	1:E:100:THR:HG22	2.48	0.48
1:E:245:PHE:O	1:E:283:THR:HG23	2.14	0.48
1:E:157:ARG:O	1:E:161:GLN:HB2	2.13	0.48
1:E:170:ILE:HG13	3:E:549:HOH:O	2.13	0.47
1:E:220:LYS:CG	2:P:4:PRO:HG3	2.44	0.47
1:E:271:PRO:HA	1:E:274:TYR:CZ	2.49	0.47
1:E:41:PRO:HA	1:E:104:SER:OG	2.15	0.47
1:E:131:GLU:O	1:E:132:TYR:HB2	2.15	0.47
1:E:262:ILE:HB	1:E:267:TYR:CE2	2.49	0.47
1:E:176:THR:HG23	1:E:324:LYS:HE2	1.96	0.47
1:E:274:TYR:HB3	3:E:491:HOH:O	2.14	0.47
1:E:51:LYS:HB3	1:E:51:LYS:HE2	1.76	0.47
1:E:33:THR:HA	1:E:123:MET:HB2	1.96	0.47
1:E:291:HIS:O	1:E:292:SER:CB	2.63	0.47
1:E:277:GLN:HE21	1:E:280:GLY:N	2.11	0.47
1:E:279:GLN:O	1:E:279:GLN:CG	2.62	0.46
1:E:99:GLN:NE2	1:E:136:VAL:HA	2.29	0.46
1:E:220:LYS:HZ2	1:E:220:LYS:HG3	1.65	0.46
1:E:274:TYR:CD1	1:E:274:TYR:O	2.68	0.46
1:E:172:PRO:HA	1:E:175:TYR:CE2	2.51	0.46
1:E:67:LEU:HD21	1:E:86:TYR:CE2	2.51	0.46
1:E:124:ALA:HB1	1:E:128:LEU:HD22	1.97	0.45
1:E:167:LEU:HG	3:E:557:HOH:O	2.16	0.45
1:E:195:ASP:O	1:E:196:SER:HB3	2.17	0.45
1:E:271:PRO:O	1:E:275:THR:HB	2.16	0.45
1:E:1:VAL:O	1:E:93:ASN:ND2	2.50	0.45
1:E:151:PHE:HB2	1:E:167:LEU:HD23	1.99	0.45
1:E:128:LEU:HD13	1:E:188:GLN:CA	2.47	0.45
1:E:128:LEU:HD13	1:E:188:GLN:CB	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:124:ALA:CB	1:E:128:LEU:HD22	2.47	0.44
1:E:25:GLN:NE2	1:E:56:PHE:HD1	2.15	0.44
1:E:199:ILE:O	1:E:200:SER:HB2	2.17	0.44
1:E:180:HIS:HB3	3:E:458:HOH:O	2.18	0.44
1:E:52:ASN:ND2	1:E:52:ASN:C	2.69	0.44
1:E:277:GLN:HB2	1:E:282:CYS:SG	2.58	0.44
1:E:269:LEU:HD11	1:E:305:PHE:CD1	2.53	0.44
1:E:244:GLU:HG3	1:E:245:PHE:H	1.83	0.44
1:E:232:GLN:HA	1:E:232:GLN:OE1	2.18	0.44
1:E:199:ILE:HG13	1:E:204:VAL:HG11	2.00	0.44
1:E:283:THR:HG22	1:E:284:SER:N	2.33	0.43
1:E:238:THR:O	1:E:245:PHE:HA	2.18	0.43
1:E:302:GLY:O	1:E:306:ILE:HG13	2.18	0.43
1:E:36:SER:OG	1:E:124:ALA:HB3	2.18	0.43
1:E:220:LYS:O	1:E:302:GLY:HA3	2.19	0.43
1:E:71:LEU:HD22	1:E:102:GLY:CA	2.43	0.43
1:E:22:THR:CA	1:E:23:PRO:C	2.86	0.43
1:E:131:GLU:O	1:E:132:TYR:CB	2.67	0.43
1:E:105:THR:O	1:E:106:GLN:HG2	2.18	0.43
1:E:220:LYS:HE3	2:P:4:PRO:HG3	2.01	0.43
1:E:184:VAL:HA	1:E:191:GLN:O	2.19	0.43
1:E:291:HIS:HB2	1:E:292:SER:H	1.60	0.42
1:E:28:THR:OG1	1:E:117:PHE:HA	2.19	0.42
1:E:255:MET:HB3	1:E:256:PRO:HD2	2.01	0.42
1:E:19:TYR:CD1	1:E:26:GLU:HB2	2.54	0.42
1:E:67:LEU:HB2	1:E:84:LEU:O	2.18	0.42
1:E:22:THR:O	1:E:61:SER:HA	2.19	0.42
1:E:199:ILE:HD12	1:E:234:ALA:CB	2.49	0.42
1:E:124:ALA:HB1	3:E:575:HOH:O	2.19	0.42
1:E:158:ASN:C	1:E:161:GLN:HB3	2.39	0.42
1:E:112:PHE:CE1	2:P:1:NOR:HE21	2.55	0.42
1:E:26:GLU:HG2	1:E:27:PHE:N	2.35	0.42
1:E:77:THR:HB	2:P:2:SMC:N	2.35	0.42
1:E:100:THR:HG21	1:E:134:ILE:HD12	2.01	0.42
1:E:82:GLY:C	1:E:83:ILE:HD13	2.40	0.42
1:E:83:ILE:N	1:E:83:ILE:CD1	2.80	0.42
1:E:137:PHE:O	1:E:140:MET:HB2	2.19	0.42
1:E:2:ALA:CB	1:E:94:ILE:CD1	2.98	0.41
1:E:36:SER:OG	1:E:128:LEU:HB3	2.19	0.41
1:E:112:PHE:HE1	1:E:120:ILE:HD11	1.85	0.41
1:E:239:GLN:HA	1:E:244:GLU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:129:ALA:CB	1:E:135:PRO:HD3	2.50	0.41
1:E:32:ASP:OD2	2:P:1:NOR:HC	2.20	0.41
1:E:38:PHE:CD2	1:E:101:VAL:HB	2.56	0.41
1:E:100:THR:CG2	1:E:134:ILE:HD12	2.50	0.41
1:E:151:PHE:CB	1:E:167:LEU:HD23	2.51	0.41
1:E:89:VAL:HG13	1:E:96:ASP:HB3	2.02	0.41
1:E:52:ASN:ND2	1:E:53:HIS:ND1	2.68	0.41
1:E:32:ASP:OD2	2:P:1:NOR:HB1	2.21	0.41
1:E:67:LEU:HD13	1:E:67:LEU:HA	1.88	0.41
1:E:20:LEU:HA	1:E:20:LEU:HD12	1.83	0.41
1:E:269:LEU:O	1:E:274:TYR:HD2	2.04	0.40
1:E:217:GLY:HA3	2:P:1:NOR:HB2	2.02	0.40
1:E:13:GLN:HE21	2:P:3:PHI:CE1	2.34	0.40
1:E:2:ALA:CB	1:E:94:ILE:HD12	2.51	0.40
1:E:123:MET:HB3	1:E:151:PHE:CZ	2.56	0.40
1:E:29:VAL:HA	1:E:119:GLY:O	2.21	0.40
1:E:1:VAL:HG11	1:E:169:ALA:HB3	2.03	0.40
1:E:47:SER:O	1:E:50:CYS:HB2	2.22	0.40
1:E:67:LEU:HD22	1:E:86:TYR:CE2	2.56	0.40
1:E:124:ALA:HA	3:E:575:HOH:O	2.22	0.40
1:E:112:PHE:HB3	3:E:399:HOH:O	2.21	0.40
1:E:192:PHE:HZ	1:E:301:LEU:HD12	1.86	0.40

All (10) 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)
3:E:527:HOH:O	3:E:527:HOH:O[2_655]	0.37	1.83
3:E:568:HOH:O	3:E:568:HOH:O[6_766]	0.58	1.62
1:E:242:TYR:CD2	1:E:244:GLU:CG[18_655]	1.47	0.73
1:E:241:GLN:O	1:E:246:ASP:OD2[18_655]	1.69	0.51
1:E:240:ASN:OD1	1:E:240:ASN:OD1[18_655]	1.87	0.33
3:E:450:HOH:O	3:E:450:HOH:O[10_546]	1.92	0.28
1:E:242:TYR:CD2	1:E:244:GLU:CD[18_655]	2.08	0.12
1:E:242:TYR:CE2	1:E:244:GLU:OE2[18_655]	2.09	0.11
3:E:521:HOH:O	3:E:521:HOH:O[17_555]	2.09	0.11
1:E:240:ASN:ND2	1:E:240:ASN:ND2[18_655]	2.11	0.09

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	E	321/323 (99%)	284 (88%)	25 (8%)	12 (4%)	<b>4</b> <b>2</b>

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	157	ARG
1	E	158	ASN
1	E	242	TYR
1	E	244	GLU
1	E	297	GLN
1	E	92	SER
1	E	159	GLY
1	E	243	GLY
1	E	290	ASN
1	E	292	SER
1	E	156	ASP
1	E	279	GLN

### 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	E	280/280 (100%)	240 (86%)	40 (14%)	<b>4</b> <b>4</b>
2	P	1/1 (100%)	1 (100%)	0	<b>100</b> <b>100</b>
All	All	281/281 (100%)	241 (86%)	40 (14%)	<b>4</b> <b>4</b>

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	17	LYS
1	E	22	THR
1	E	29	VAL
1	E	51	LYS
1	E	52	ASN
1	E	59	ARG
1	E	67	LEU
1	E	69	LYS
1	E	77	THR
1	E	80	MET
1	E	81	GLN
1	E	83	ILE
1	E	93	ASN
1	E	105	THR
1	E	110	ASP
1	E	116	GLU
1	E	126	PRO
1	E	128	LEU
1	E	148	GLN
1	E	157	ARG
1	E	161	GLN
1	E	162	GLU
1	E	219	SER
1	E	220	LYS
1	E	225	SER
1	E	226	SER
1	E	229	LEU
1	E	231	ILE
1	E	242	TYR
1	E	245	PHE
1	E	247	ILE
1	E	250	ASP
1	E	253	SER
1	E	272	SER
1	E	281	PHE
1	E	284	SER
1	E	292	SER
1	E	297	GLN
1	E	298	LYS
1	E	311	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such

sidechains are listed below:

Mol	Chain	Res	Type
1	E	25	GLN
1	E	52	ASN
1	E	54	GLN
1	E	81	GLN
1	E	99	GLN
1	E	139	ASN
1	E	142	ASN
1	E	144	HIS
1	E	148	GLN
1	E	161	GLN
1	E	251	ASN
1	E	277	GLN
1	E	318	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	$\# Z  > 2$
2	NOR	P	1	2	17,17,17	2.94	1 (5%)	18,22,22	2.74	5 (27%)
2	SMC	P	2	2	5,6,7	0.47	0	2,6,8	4.25	2 (100%)
2	PHI	P	3	2	11,12,13	8.47	1 (9%)	12,15,17	2.52	7 (58%)

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	NOR	P	1	2	-	0/16/24/24	0/1/1/1
2	SMC	P	2	2	-	0/3/5/7	0/0/0/0
2	PHI	P	3	2	-	0/4/6/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	3	PHI	CZ-I	-28.00	1.36	2.10
2	P	1	NOR	CH-CA	-11.70	1.41	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	1	NOR	CE2-CD2-CG	-8.31	98.82	112.22
2	P	3	PHI	CE2-CZ-CE1	-5.02	112.95	120.66
2	P	1	NOR	CM-O2-C	-4.79	111.84	117.87
2	P	3	PHI	O-C-CA	-2.40	119.23	125.49
2	P	2	SMC	O-C-CA	-2.38	119.29	125.49
2	P	3	PHI	CD2-CG-CD1	-2.35	114.37	118.13
2	P	3	PHI	CG-CB-CA	2.25	119.30	114.21
2	P	1	NOR	CD2-CG-CD1	2.57	115.71	109.26
2	P	3	PHI	CD2-CE2-CZ	2.57	123.34	119.53
2	P	1	NOR	O2-CM-CM2	2.69	114.08	107.12
2	P	3	PHI	CE1-CZ-I	3.00	124.39	119.67
2	P	3	PHI	CD1-CE1-CZ	3.28	124.40	119.53
2	P	1	NOR	O2-C-CH	3.93	118.64	111.14
2	P	2	SMC	CS-SG-CB	5.52	110.18	101.21

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	1	NOR	6	0
2	P	2	SMC	1	0
2	P	3	PHI	8	0

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	E	323/323 (100%)	0.94	28 (8%) 13 18	0, 0, 0, 1	0
2	P	1/4 (25%)	7.61	1 (100%) 0 0	1, 1, 1, 1	0
All	All	324/327 (99%)	0.96	29 (8%) 12 17	0, 0, 0, 1	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	242	TYR	11.4
1	E	241	GLN	7.8
2	P	4	PRO	7.6
1	E	159	GLY	7.0
1	E	290	ASN	6.5
1	E	292	SER	5.8
1	E	291	HIS	5.8
1	E	161	GLN	5.8
1	E	156	ASP	4.6
1	E	158	ASN	4.6
1	E	243	GLY	4.5
1	E	244	GLU	4.2
1	E	132	TYR	3.6
1	E	157	ARG	3.2
1	E	240	ASN	3.1
1	E	281	PHE	2.9
1	E	239	GLN	2.8
1	E	9	TYR	2.7
1	E	250	ASP	2.6
1	E	39	TRP	2.4
1	E	229	LEU	2.3
1	E	297	GLN	2.3
1	E	102	GLY	2.3
1	E	207	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	153	VAL	2.3
1	E	150	LEU	2.2
1	E	41	PRO	2.2
1	E	106	GLN	2.1
1	E	238	THR	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	PHI	P	3	12/13	0.96	0.19	-	0,0,0,1	0
2	NOR	P	1	17/17	0.83	0.21	-	0,0,0,0	0
2	SMC	P	2	7/8	0.92	0.14	-	0,0,0,0	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.