



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:07 AM GMT

PDB ID : 1A0O
Title : CHEY-BINDING DOMAIN OF CHEA IN COMPLEX WITH CHEY
Authors : Chinardet, N.; Welch, M.; Mourey, L.; Birck, C.; Samama, J.P.
Deposited on : 1997-12-05
Resolution : 2.95 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

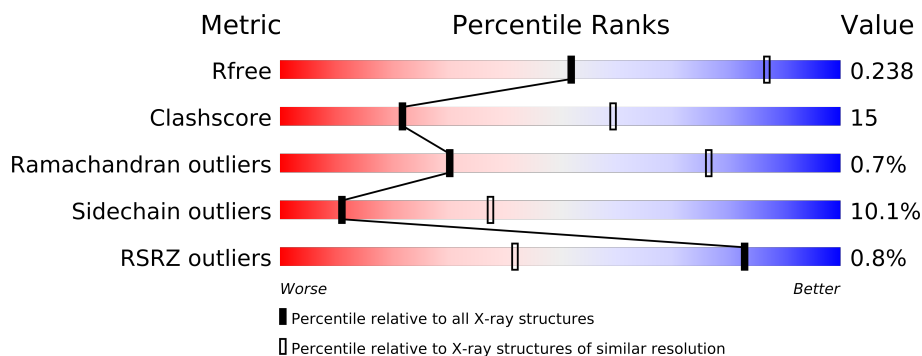
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.95 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}	66092	1587 (3.00-2.92)
Clashscore	79885	2029 (3.00-2.92)
Ramachandran outliers	78287	1955 (3.00-2.92)
Sidechain outliers	78261	1958 (3.00-2.92)
RSRZ outliers	66119	1588 (3.00-2.92)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	128	
1	C	128	
1	E	128	
1	G	128	
2	B	134	
2	D	134	
2	F	134	
2	H	134	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5969 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CHEY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	128	Total	C	N	O	S	0	0	0
			979	623	162	188	6			
1	C	128	Total	C	N	O	S	0	0	0
			979	623	162	188	6			
1	E	128	Total	C	N	O	S	0	0	0
			979	623	162	188	6			
1	G	128	Total	C	N	O	S	0	0	0
			979	623	162	188	6			

- Molecule 2 is a protein called CHEA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	70	Total	C	N	O	S	0	0	0
			524	329	84	110	1			
2	D	67	Total	C	N	O	S	0	0	0
			505	318	81	105	1			
2	F	69	Total	C	N	O	S	0	0	0
			515	324	83	107	1			
2	H	67	Total	C	N	O	S	0	0	0
			505	318	81	105	1			

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Mn	0	0
			1	1		
3	A	1	Total	Mn	0	0
			1	1		
3	C	1	Total	Mn	0	0
			1	1		
3	E	1	Total	Mn	0	0
			1	1		

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.

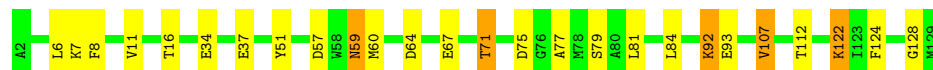
- Molecule 1: CHEY

Chain A: 



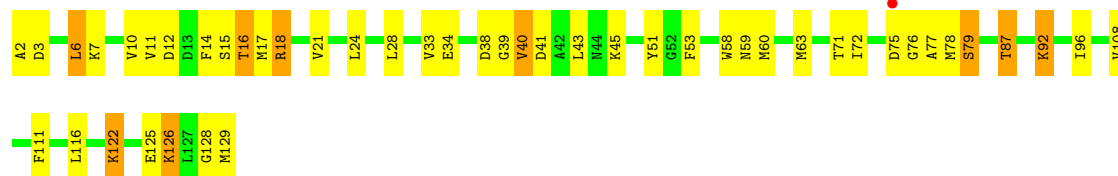
- Molecule 1: CHEY

Chain C: 



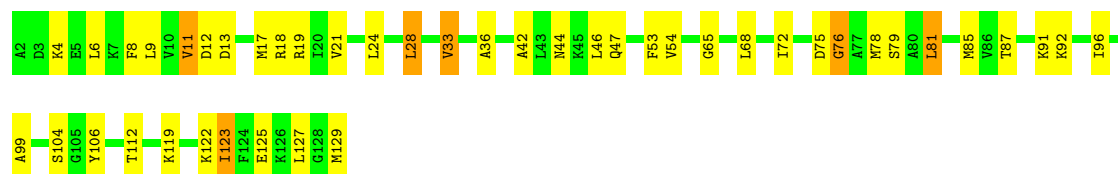
- Molecule 1: CHEY

Chain E: 



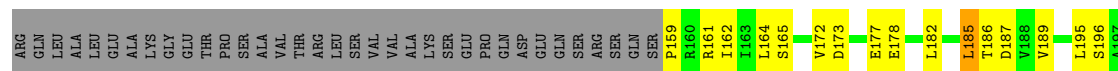
- Molecule 1: CHEY

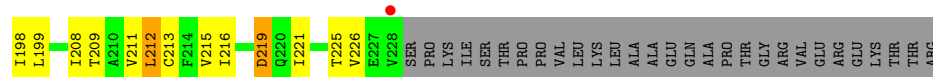
Chain G: 



- Molecule 2: CHEA

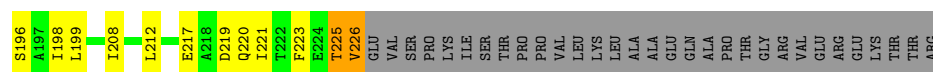
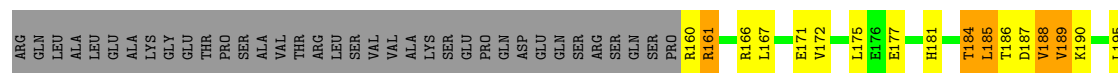
Chain B: 





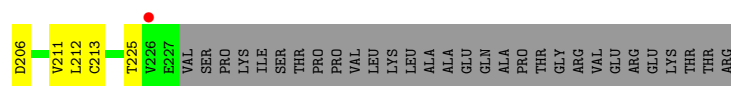
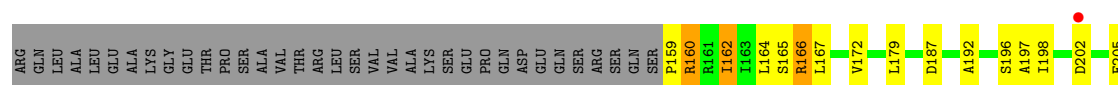
• Molecule 2: CHEA

Chain D:



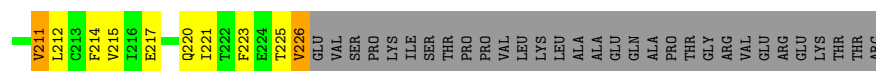
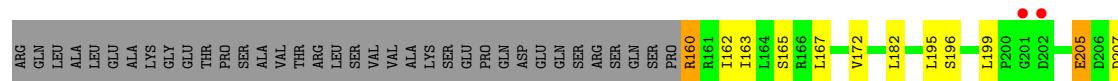
• Molecule 2: CHEA

Chain F:



• Molecule 2: CHEA

Chain H:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.90Å 156.97Å 65.97Å 90.00° 91.70° 90.00°	Depositor
Resolution (Å)	27.00 – 2.95 26.79 – 2.95	Depositor EDS
% Data completeness (in resolution range)	90.9 (27.00-2.95) 91.0 (26.79-2.95)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.08 (at 2.94Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.187 , 0.235 0.195 , 0.238	Depositor DCC
R_{free} test set	1073 reflections (5.11%)	DCC
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.336	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 22.4	EDS
Estimated twinning fraction	0.207 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 20984 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5969	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.39	0/991	0.59	0/1332
1	C	0.40	0/991	0.60	0/1332
1	E	0.39	0/991	0.59	0/1332
1	G	0.38	0/991	0.58	0/1332
2	B	0.37	0/527	0.63	1/716 (0.1%)
2	D	0.40	0/508	0.63	0/690
2	F	0.36	0/518	0.68	1/704 (0.1%)
2	H	0.41	0/508	0.66	0/690
All	All	0.39	0/6025	0.61	2/8128 (0.0%)

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	A	0	1
1	E	0	1
1	G	0	1
2	D	0	3
2	F	0	2
2	H	0	1
All	All	0	9

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	159	PRO	N-CA-CB	5.69	110.13	103.30
2	F	159	PRO	N-CA-CB	5.54	109.95	103.30

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	19	ARG	Sidechain
2	D	160	ARG	Sidechain
2	D	161	ARG	Sidechain
2	D	166	ARG	Sidechain
1	E	18	ARG	Sidechain
2	F	160	ARG	Sidechain
2	F	166	ARG	Sidechain
1	G	19	ARG	Sidechain
2	H	160	ARG	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	979	0	999	22	0
1	C	979	0	999	19	0
1	E	979	0	999	37	0
1	G	979	0	999	38	0
2	B	524	0	520	16	0
2	D	505	0	510	27	0
2	F	515	0	514	11	0
2	H	505	0	510	17	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
All	All	5969	0	6050	182	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 15.

All (182) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:225:THR:O	2:D:226:VAL:HG12	1.28	1.23
1:E:92:LYS:HD3	1:E:96:ILE:HD11	1.26	1.17
1:E:92:LYS:HD3	1:E:96:ILE:CD1	1.79	1.11
2:D:225:THR:O	2:D:226:VAL:CG1	2.01	1.08
1:C:84:LEU:HD11	1:C:107:VAL:HG12	1.29	1.08
1:G:106:TYR:O	1:G:119:LYS:HE3	1.64	0.97
2:D:185:LEU:N	2:D:185:LEU:HD22	1.79	0.96
2:F:162:ILE:HD13	2:F:164:LEU:HD21	1.44	0.96
1:E:28:LEU:CD1	1:E:116:LEU:HD23	2.06	0.86
1:E:28:LEU:HD11	1:E:116:LEU:HD23	1.56	0.85
1:A:72:ILE:HG12	1:A:78:MET:CE	2.07	0.85
1:C:84:LEU:HD11	1:C:107:VAL:CG1	2.09	0.82
2:F:162:ILE:CD1	2:F:164:LEU:HD21	2.09	0.82
2:D:185:LEU:N	2:D:185:LEU:CD2	2.43	0.81
2:H:182:LEU:HD11	2:H:211:VAL:HG21	1.64	0.80
1:C:7:LYS:HE2	1:C:34:GLU:HG3	1.64	0.78
1:A:72:ILE:HA	1:A:78:MET:CE	2.12	0.78
1:A:6:LEU:HD11	1:A:52:GLY:HA3	1.66	0.78
2:D:190:LYS:HG2	2:D:195:LEU:HD12	1.64	0.77
2:H:160:ARG:NH1	2:H:205:GLU:OE1	2.18	0.76
2:D:186:THR:HB	2:D:198:ILE:HG13	1.68	0.76
1:A:72:ILE:HG12	1:A:78:MET:HE3	1.68	0.75
1:G:75:ASP:OD2	1:G:78:MET:HG2	1.87	0.75
2:D:225:THR:C	2:D:226:VAL:CG1	2.51	0.74
1:G:17:MET:O	1:G:21:VAL:HG23	1.88	0.74
1:G:68:LEU:O	1:G:72:ILE:HG13	1.88	0.74
1:C:84:LEU:CD1	1:C:107:VAL:HG12	2.15	0.72
1:E:24:LEU:O	1:E:28:LEU:HG	1.90	0.71
1:E:92:LYS:HD3	1:E:96:ILE:HD12	1.68	0.71
1:E:92:LYS:CD	1:E:96:ILE:CD1	2.67	0.70
1:E:58:TRP:HD1	1:E:59:ASN:OD1	1.74	0.70
2:F:162:ILE:HD13	2:F:164:LEU:CD2	2.21	0.68
1:A:72:ILE:HG12	1:A:78:MET:HE1	1.73	0.67
2:H:217:GLU:H	2:H:220:GLN:NE2	1.93	0.66
1:E:34:GLU:OE1	1:E:51:TYR:OH	2.10	0.66
2:D:208:ILE:O	2:D:212:LEU:HG	1.96	0.66
1:C:59:ASN:OD1	1:C:59:ASN:N	2.22	0.65
2:D:186:THR:HG22	2:D:187:ASP:N	2.11	0.65
1:G:53:PHE:CE2	1:G:123:ILE:HG21	2.31	0.65
2:D:187:ASP:H	2:D:198:ILE:HG12	1.61	0.65
2:D:171:GLU:O	2:D:175:LEU:HD12	1.97	0.65
2:B:172:VAL:HG23	2:B:195:LEU:HB2	1.79	0.64
1:E:2:ALA:N	1:E:129:MET:OXT	2.31	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:64:ASP:OD1	1:C:67:GLU:HG3	1.99	0.63
2:D:190:LYS:HG2	2:D:195:LEU:CD1	2.29	0.63
1:G:9:LEU:HD23	1:G:46:LEU:HD21	1.79	0.62
2:B:162:ILE:HB	2:B:199:LEU:HD11	1.80	0.62
2:D:185:LEU:H	2:D:185:LEU:CD2	2.12	0.62
1:G:42:ALA:O	1:G:46:LEU:HG	1.99	0.62
2:H:182:LEU:HD23	2:H:182:LEU:N	2.14	0.62
1:A:7:LYS:HB3	1:A:51:TYR:HA	1.82	0.62
1:C:6:LEU:HD23	1:C:8:PHE:CE1	2.34	0.61
2:B:185:LEU:HD12	2:B:185:LEU:H	1.65	0.61
2:H:182:LEU:HD11	2:H:211:VAL:CG2	2.31	0.60
1:E:17:MET:O	1:E:21:VAL:HG23	2.00	0.60
2:B:164:LEU:HD12	2:B:195:LEU:HD23	1.83	0.60
1:E:92:LYS:CD	1:E:96:ILE:HD12	2.29	0.60
1:G:127:LEU:HD13	1:G:129:MET:HE1	1.84	0.59
1:G:75:ASP:OD1	1:G:76:GLY:N	2.35	0.59
1:E:40:VAL:HG12	1:E:41:ASP:N	2.18	0.59
2:D:219:ASP:OD1	2:D:219:ASP:C	2.42	0.58
1:G:127:LEU:HD13	1:G:129:MET:CE	2.33	0.58
1:G:53:PHE:HE2	1:G:123:ILE:HG21	1.68	0.58
2:D:225:THR:C	2:D:226:VAL:HG13	2.23	0.57
2:H:160:ARG:NH1	2:H:205:GLU:CD	2.58	0.57
1:E:75:ASP:O	1:E:79:SER:HB2	2.05	0.57
1:A:72:ILE:HA	1:A:78:MET:HE2	1.85	0.57
1:E:28:LEU:HD12	1:E:116:LEU:HD23	1.83	0.57
1:G:54:VAL:HG23	1:G:81:LEU:HD12	1.85	0.57
2:B:208:ILE:HG22	2:B:212:LEU:HD22	1.85	0.57
1:G:8:PHE:HB2	1:G:33:VAL:HG13	1.86	0.56
1:E:75:ASP:OD2	1:E:78:MET:CE	2.53	0.56
1:C:7:LYS:HE2	1:C:34:GLU:CG	2.36	0.55
2:D:221:ILE:HD11	2:D:223:PHE:CZ	2.42	0.55
1:G:9:LEU:CD2	1:G:46:LEU:HD21	2.37	0.55
1:G:4:LYS:NZ	1:G:28:LEU:O	2.39	0.55
2:F:179:LEU:HD21	2:F:212:LEU:HD11	1.88	0.55
1:A:6:LEU:HD21	1:A:53:PHE:HB2	1.89	0.54
1:G:75:ASP:O	1:G:79:SER:HB3	2.07	0.54
1:A:53:PHE:HA	1:A:82:PRO:HG2	1.88	0.54
1:G:122:LYS:NZ	1:G:125:GLU:OE2	2.40	0.54
1:G:6:LEU:HD21	1:G:53:PHE:HB2	1.89	0.54
2:B:182:LEU:HD21	2:B:211:VAL:HG21	1.90	0.54
2:D:184:THR:C	2:D:185:LEU:HD22	2.28	0.54
1:C:67:GLU:O	1:C:71:THR:OG1	2.26	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:75:ASP:OD2	1:E:78:MET:HE3	2.07	0.53
1:G:127:LEU:CD1	1:G:129:MET:HE1	2.38	0.53
2:B:161:ARG:HB2	2:B:198:ILE:HD13	1.90	0.53
1:G:127:LEU:CD1	1:G:129:MET:CE	2.87	0.53
1:A:72:ILE:HG23	1:A:78:MET:HE3	1.90	0.52
1:E:14:PHE:HD2	1:E:17:MET:HB2	1.75	0.52
1:C:122:LYS:HE2	1:C:122:LYS:HA	1.90	0.52
1:A:6:LEU:CD1	1:A:52:GLY:HA3	2.36	0.52
1:E:11:VAL:HG12	1:E:60:MET:SD	2.50	0.52
2:B:209:THR:HG23	2:B:221:ILE:HD12	1.92	0.52
2:F:167:LEU:HD12	2:F:172:VAL:HG22	1.92	0.52
1:A:6:LEU:HD11	1:A:52:GLY:CA	2.38	0.51
2:B:162:ILE:O	2:B:196:SER:HA	2.10	0.51
2:F:187:ASP:HB2	2:F:198:ILE:HD12	1.93	0.51
1:E:12:ASP:O	1:E:18:ARG:HD3	2.11	0.50
1:A:126:LYS:O	1:A:127:LEU:HD23	2.11	0.50
2:H:207:ASP:O	2:H:211:VAL:HG13	2.11	0.50
2:D:167:LEU:HD12	2:D:172:VAL:HG22	1.94	0.50
1:A:85:MET:O	1:A:106:TYR:HA	2.12	0.49
1:A:89:GLU:HG2	1:A:91:LYS:HG2	1.94	0.49
1:A:89:GLU:CG	1:A:91:LYS:HG2	2.41	0.49
1:A:122:LYS:HD3	1:A:122:LYS:O	2.12	0.49
2:B:162:ILE:HB	2:B:199:LEU:CD1	2.42	0.49
2:H:225:THR:O	2:H:226:VAL:HG12	2.11	0.49
1:C:75:ASP:OD1	1:C:77:ALA:N	2.46	0.49
2:B:185:LEU:N	2:B:185:LEU:HD12	2.28	0.48
2:B:213:CYS:HA	2:B:216:ILE:O	2.13	0.48
2:H:167:LEU:HD12	2:H:172:VAL:HG22	1.94	0.48
1:C:7:LYS:HG2	1:C:51:TYR:CD2	2.49	0.48
2:H:217:GLU:HB2	2:H:220:GLN:HE21	1.78	0.48
2:H:162:ILE:HD11	2:H:221:ILE:HG23	1.94	0.47
1:G:12:ASP:HB3	1:G:18:ARG:HG2	1.96	0.47
1:G:99:ALA:HB2	2:H:214:PHE:CZ	2.49	0.47
1:C:11:VAL:HG12	1:C:60:MET:SD	2.54	0.47
1:A:43:LEU:HD21	1:A:71:THR:HG21	1.97	0.47
1:E:111:PHE:CE1	1:E:116:LEU:HD13	2.50	0.47
1:E:38:ASP:HA	1:E:60:MET:SD	2.54	0.47
1:E:122:LYS:HA	1:E:122:LYS:HD3	1.44	0.46
1:A:122:LYS:HE3	2:B:215:VAL:HG13	1.96	0.46
1:E:3:ASP:O	1:E:6:LEU:HB2	2.16	0.46
1:G:9:LEU:CD2	1:G:46:LEU:CD2	2.94	0.46
1:A:72:ILE:CG1	1:A:78:MET:CE	2.89	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:28:LEU:HD12	1:G:28:LEU:HA	1.75	0.46
1:C:75:ASP:O	1:C:79:SER:HB3	2.15	0.46
2:D:219:ASP:OD1	2:D:220:GLN:N	2.50	0.45
2:F:187:ASP:O	2:F:197:ALA:HA	2.17	0.45
2:H:160:ARG:HH12	2:H:205:GLU:CD	2.18	0.45
2:F:160:ARG:HD3	2:F:205:GLU:OE2	2.17	0.45
1:C:92:LYS:HB3	2:D:181:HIS:O	2.16	0.45
1:G:65:GLY:HA3	1:G:85:MET:HE3	1.99	0.45
1:E:14:PHE:CE2	1:E:16:THR:HB	2.52	0.44
2:F:162:ILE:O	2:F:196:SER:HA	2.18	0.44
1:G:68:LEU:HG	1:G:72:ILE:HD11	2.00	0.44
1:G:9:LEU:HD22	1:G:46:LEU:CD2	2.48	0.44
1:E:76:GLY:O	1:E:77:ALA:HB3	2.17	0.44
1:A:2:ALA:N	1:A:129:MET:O	2.51	0.44
1:E:126:LYS:NZ	2:F:213:CYS:O	2.47	0.43
1:G:4:LYS:HD3	1:G:4:LYS:HA	1.70	0.43
1:C:122:LYS:HE3	1:C:122:LYS:HB2	1.84	0.43
2:D:217:GLU:H	2:D:220:GLN:NE2	2.16	0.43
1:C:75:ASP:OD1	1:C:75:ASP:C	2.57	0.43
2:H:162:ILE:O	2:H:196:SER:HA	2.18	0.43
1:G:11:VAL:HG13	1:G:36:ALA:HB3	2.00	0.43
1:C:124:PHE:O	1:C:128:GLY:N	2.51	0.42
2:D:189:VAL:HG23	2:D:196:SER:OG	2.20	0.42
1:E:43:LEU:HD21	1:E:71:THR:HG21	2.01	0.42
1:G:92:LYS:O	1:G:96:ILE:HG12	2.19	0.42
2:H:205:GLU:HG3	2:H:223:PHE:CE2	2.55	0.42
2:H:212:LEU:HD23	2:H:212:LEU:HA	1.79	0.42
2:D:161:ARG:HH11	2:D:226:VAL:HG11	1.83	0.42
2:D:189:VAL:CG2	2:D:196:SER:OG	2.68	0.42
2:B:177:GLU:HG3	2:B:178:GLU:N	2.35	0.42
1:G:12:ASP:HB3	1:G:18:ARG:CG	2.50	0.41
1:E:58:TRP:CE3	1:E:87:THR:HG23	2.55	0.41
1:G:122:LYS:HE3	2:H:215:VAL:O	2.21	0.41
2:B:219:ASP:OD1	2:B:219:ASP:C	2.58	0.41
1:E:6:LEU:CD1	1:E:53:PHE:HB2	2.51	0.41
1:E:6:LEU:HD11	1:E:53:PHE:HB2	2.02	0.41
1:G:127:LEU:CB	1:G:129:MET:HE3	2.51	0.41
1:E:7:LYS:HB3	1:E:51:TYR:HA	2.02	0.41
1:G:127:LEU:CD1	1:G:129:MET:HE3	2.51	0.41
1:E:122:LYS:HE3	1:E:125:GLU:OE1	2.21	0.41
2:D:186:THR:O	2:D:188:VAL:HG23	2.21	0.41
2:D:198:ILE:O	2:D:198:ILE:HG13	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:44:ASN:O	1:G:47:GLN:HB2	2.20	0.41
2:F:160:ARG:NH1	2:F:205:GLU:OE2	2.44	0.41
1:E:126:LYS:HB3	1:E:126:LYS:HE2	1.85	0.41
1:E:72:ILE:HG23	1:E:78:MET:HB3	2.03	0.40
1:G:12:ASP:O	1:G:18:ARG:HD3	2.22	0.40
1:E:10:VAL:CG2	1:E:33:VAL:CG1	2.99	0.40
2:B:172:VAL:HG13	2:B:173:ASP:N	2.36	0.40
1:G:92:LYS:HB3	1:G:92:LYS:HE2	1.86	0.40
1:E:39:GLY:HA3	1:E:63:MET:HB3	2.03	0.40
1:G:13:ASP:OD1	1:G:13:ASP:N	2.52	0.40
1:A:11:VAL:HG12	1:A:60:MET:SD	2.61	0.40
2:D:187:ASP:H	2:D:198:ILE:CG1	2.32	0.40
1:C:93:GLU:H	1:C:93:GLU:CD	2.24	0.40

There are no symmetry-related clashes.

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	A	126/128 (98%)	117 (93%)	9 (7%)	0	100	100
1	C	126/128 (98%)	120 (95%)	6 (5%)	0	100	100
1	E	126/128 (98%)	117 (93%)	7 (6%)	2 (2%)	14	54
1	G	126/128 (98%)	120 (95%)	5 (4%)	1 (1%)	27	74
2	B	68/134 (51%)	64 (94%)	4 (6%)	0	100	100
2	D	65/134 (48%)	59 (91%)	6 (9%)	0	100	100
2	F	67/134 (50%)	59 (88%)	6 (9%)	2 (3%)	7	32
2	H	65/134 (48%)	57 (88%)	8 (12%)	0	100	100
All	All	769/1048 (73%)	713 (93%)	51 (7%)	5 (1%)	30	76

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	79	SER
2	F	192	ALA
1	E	128	GLY
1	G	76	GLY
2	F	202	ASP

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	102/102 (100%)	98 (96%)	4 (4%)	43	83
1	C	102/102 (100%)	92 (90%)	10 (10%)	12	39
1	E	102/102 (100%)	92 (90%)	10 (10%)	12	39
1	G	102/102 (100%)	92 (90%)	10 (10%)	12	39
2	B	57/115 (50%)	48 (84%)	9 (16%)	4	16
2	D	56/115 (49%)	48 (86%)	8 (14%)	5	20
2	F	56/115 (49%)	50 (89%)	6 (11%)	10	33
2	H	56/115 (49%)	49 (88%)	7 (12%)	7	26
All	All	633/868 (73%)	569 (90%)	64 (10%)	11	37

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	LEU
1	A	47	GLN
1	A	59	ASN
1	A	91	LYS
2	B	165	SER
2	B	185	LEU
2	B	186	THR
2	B	187	ASP
2	B	189	VAL
2	B	212	LEU
2	B	219	ASP
2	B	225	THR
2	B	226	VAL

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Mol	Chain	Res	Type
1	C	16	THR
1	C	37	GLU
1	C	57	ASP
1	C	59	ASN
1	C	71	THR
1	C	81	LEU
1	C	92	LYS
1	C	107	VAL
1	C	112	THR
1	C	122	LYS
2	D	177	GLU
2	D	184	THR
2	D	185	LEU
2	D	188	VAL
2	D	189	VAL
2	D	199	LEU
2	D	225	THR
2	D	226	VAL
1	E	6	LEU
1	E	15	SER
1	E	16	THR
1	E	40	VAL
1	E	45	LYS
1	E	87	THR
1	E	92	LYS
1	E	108	VAL
1	E	122	LYS
1	E	126	LYS
2	F	162	ILE
2	F	165	SER
2	F	166	ARG
2	F	206	ASP
2	F	211	VAL
2	F	225	THR
1	G	11	VAL
1	G	24	LEU
1	G	28	LEU
1	G	33	VAL
1	G	81	LEU
1	G	87	THR
1	G	91	LYS
1	G	104	SER

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Mol	Chain	Res	Type
1	G	112	THR
1	G	123	ILE
2	H	163	ILE
2	H	165	SER
2	H	195	LEU
2	H	199	LEU
2	H	205	GLU
2	H	211	VAL
2	H	226	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	32	ASN
1	A	94	ASN
2	B	220	GLN
1	C	47	GLN
2	D	220	GLN
1	E	121	ASN
2	H	220	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	128/128 (100%)	-0.32	0 100 100	8, 23, 53, 88	0
1	C	128/128 (100%)	-0.36	0 100 100	8, 24, 53, 68	0
1	E	128/128 (100%)	-0.21	1 (0%) 83 42	8, 24, 60, 80	0
1	G	128/128 (100%)	-0.29	0 100 100	9, 22, 59, 82	0
2	B	70/134 (52%)	0.07	1 (1%) 72 34	17, 40, 81, 117	0
2	D	67/134 (50%)	-0.20	0 100 100	15, 38, 65, 101	0
2	F	69/134 (51%)	-0.07	2 (2%) 49 22	16, 44, 82, 107	0
2	H	67/134 (50%)	-0.20	2 (2%) 48 21	13, 38, 74, 95	0
All	All	785/1048 (74%)	-0.23	6 (0%) 83 42	8, 28, 68, 117	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	228	VAL	4.7
2	H	201	GLY	2.8
2	H	202	ASP	2.3
1	E	75	ASP	2.1
2	F	202	ASP	2.0
2	F	226	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MN	E	130	1/1	0.12	-2.53	25,25,25,25	0
3	MN	C	1	1/1	0.08	-2.81	28,28,28,28	0
3	MN	G	130	1/1	0.11	-2.92	27,27,27,27	0
3	MN	A	130	1/1	0.13	-3.23	28,28,28,28	0

6.5 Other polymers

There are no such residues in this entry.