Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2023 – 07:22 AM EDT

PDB ID	:	3BCN
Title	:	Crystal structure of a papain-like cysteine protease Ervatamin-A complexed with irreversible inhibitor E-64
Authors	:	Ghosh, R.; Chakrabarti, C.; Dattagupta, J.K.; Biswas, S.
Deposited on	:	2007-11-13
Resolution	:	2.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

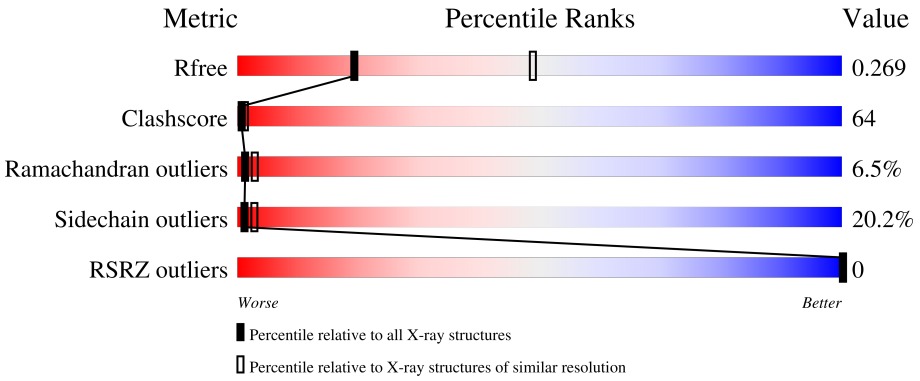
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	209	25% 56% 18%
1	B	209	30% 57% 12%

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	E64	B	215	X	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3302 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 Ervatamin-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	0	0
			1604	1012	293	288	11			
1	B	209	Total	C	N	O	S	0	0	0
			1604	1012	293	288	11			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	LEU	-	SEE REMARK 999	UNP A5YVK8
A	2	PRO	-	SEE REMARK 999	UNP A5YVK8
A	3	GLU	-	SEE REMARK 999	UNP A5YVK8
A	4	HIS	-	SEE REMARK 999	UNP A5YVK8
A	5	VAL	-	SEE REMARK 999	UNP A5YVK8
A	6	ASP	-	SEE REMARK 999	UNP A5YVK8
A	7	TRP	-	SEE REMARK 999	UNP A5YVK8
A	8	ARG	-	SEE REMARK 999	UNP A5YVK8
A	9	ALA	-	SEE REMARK 999	UNP A5YVK8
A	10	LYS	-	SEE REMARK 999	UNP A5YVK8
A	11	GLY	-	SEE REMARK 999	UNP A5YVK8
A	144	GLY	SER	SEE REMARK 999	UNP A5YVK8
A	196	CYS	-	SEE REMARK 999	UNP A5YVK8
A	197	GLY	-	SEE REMARK 999	UNP A5YVK8
A	198	ILE	-	SEE REMARK 999	UNP A5YVK8
A	199	ALA	-	SEE REMARK 999	UNP A5YVK8
A	200	ARG	-	SEE REMARK 999	UNP A5YVK8
A	201	LEU	-	SEE REMARK 999	UNP A5YVK8
A	202	PRO	-	SEE REMARK 999	UNP A5YVK8
A	203	PHE	-	SEE REMARK 999	UNP A5YVK8
A	204	TYR	-	SEE REMARK 999	UNP A5YVK8
A	205	PRO	-	SEE REMARK 999	UNP A5YVK8
A	206	THR	-	SEE REMARK 999	UNP A5YVK8
A	207	LYS	-	SEE REMARK 999	UNP A5YVK8
A	208	ALA	-	SEE REMARK 999	UNP A5YVK8

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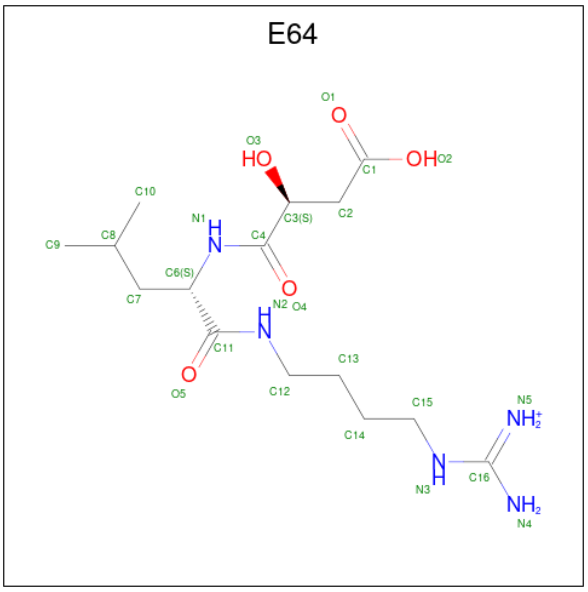
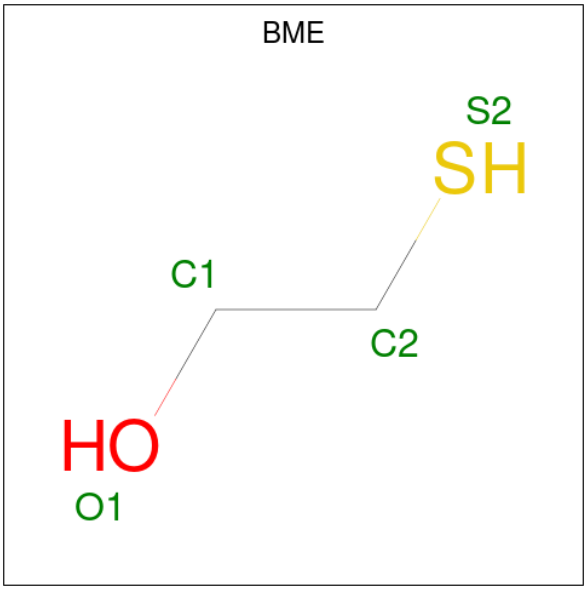
Chain	Residue	Modelled	Actual	Comment	Reference
A	209	UNK	-	SEE REMARK 999	UNP A5YVK8
B	1	LEU	-	SEE REMARK 999	UNP A5YVK8
B	2	PRO	-	SEE REMARK 999	UNP A5YVK8
B	3	GLU	-	SEE REMARK 999	UNP A5YVK8
B	4	HIS	-	SEE REMARK 999	UNP A5YVK8
B	5	VAL	-	SEE REMARK 999	UNP A5YVK8
B	6	ASP	-	SEE REMARK 999	UNP A5YVK8
B	7	TRP	-	SEE REMARK 999	UNP A5YVK8
B	8	ARG	-	SEE REMARK 999	UNP A5YVK8
B	9	ALA	-	SEE REMARK 999	UNP A5YVK8
B	10	LYS	-	SEE REMARK 999	UNP A5YVK8
B	11	GLY	-	SEE REMARK 999	UNP A5YVK8
B	144	GLY	SER	SEE REMARK 999	UNP A5YVK8
B	196	CYS	-	SEE REMARK 999	UNP A5YVK8
B	197	GLY	-	SEE REMARK 999	UNP A5YVK8
B	198	ILE	-	SEE REMARK 999	UNP A5YVK8
B	199	ALA	-	SEE REMARK 999	UNP A5YVK8
B	200	ARG	-	SEE REMARK 999	UNP A5YVK8
B	201	LEU	-	SEE REMARK 999	UNP A5YVK8
B	202	PRO	-	SEE REMARK 999	UNP A5YVK8
B	203	PHE	-	SEE REMARK 999	UNP A5YVK8
B	204	TYR	-	SEE REMARK 999	UNP A5YVK8
B	205	PRO	-	SEE REMARK 999	UNP A5YVK8
B	206	THR	-	SEE REMARK 999	UNP A5YVK8
B	207	LYS	-	SEE REMARK 999	UNP A5YVK8
B	208	ALA	-	SEE REMARK 999	UNP A5YVK8
B	209	UNK	-	SEE REMARK 999	UNP A5YVK8

- Molecule 2 is N-[N-[1-HYDROXYCARBOXYETHYL-CARBONYL]LEUCYLAMINO-BUTYL]-GUANIDINE (three-letter code: E64) (formula: $C_{15}H_{30}N_5O_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	25	15	5	5	0	0
			Total	C	N	O		
2	B	1	25	15	5	5	0	0

- Molecule 3 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C_2H_6OS).



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

- Molecule 4 is water.

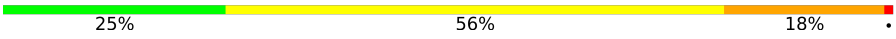
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	23	Total O 23 23	0	0
4	B	17	Total O 17 17	0	0

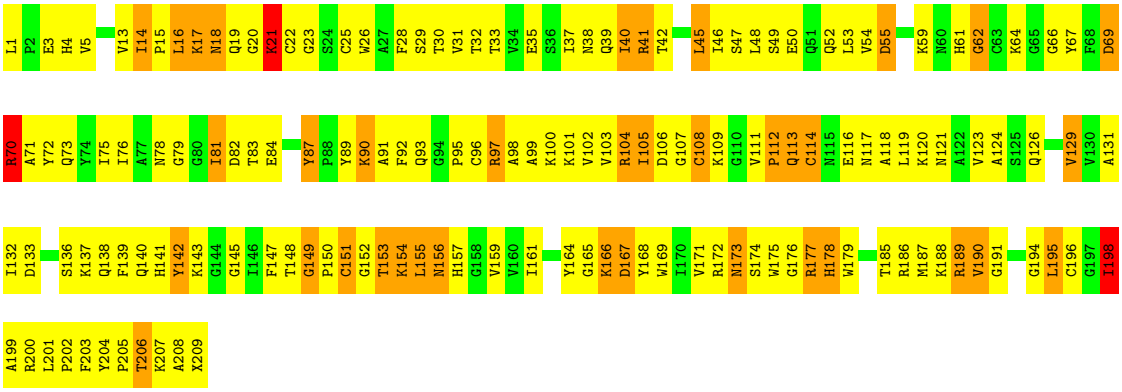
3 Residue-property plots (i)

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: Ervatamin-A

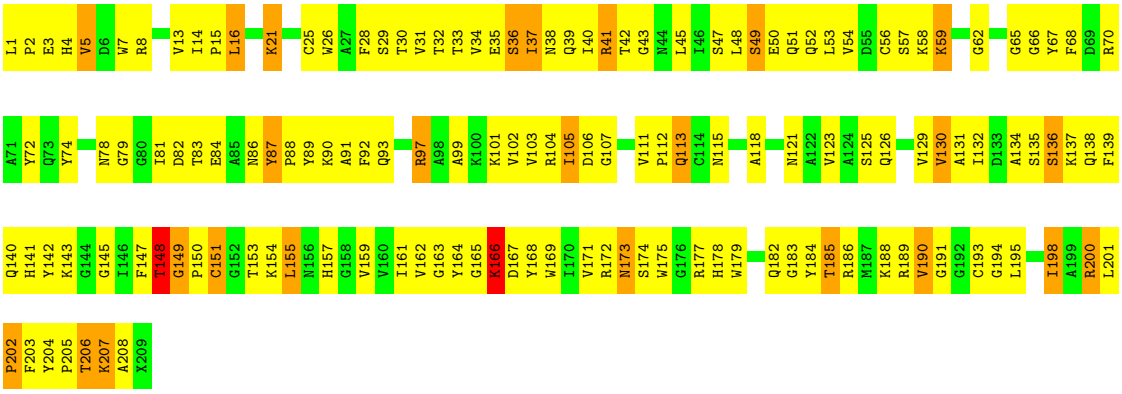
Chain A:





• Molecule 1: Ervatamin-A

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	31.17Å 105.59Å 73.93Å 90.00° 101.96° 90.00°	Depositor
Resolution (Å)	30.00 – 2.85 42.64 – 2.50	Depositor EDS
% Data completeness (in resolution range)	75.0 (30.00-2.85) 85.1 (42.64-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.31 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.240 , 0.270 0.245 , 0.269	Depositor DCC
R_{free} test set	708 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	26.3	Xtriage
Anisotropy	0.764	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 5.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.427 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.85	EDS
Total number of atoms	3302	wwPDB-VP
Average B, all atoms (Å^2)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.22% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: E64, BME

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	0.48	0/1637	0.71	1/2212 (0.0%)
1	B	0.46	0/1637	0.71	0/2212
All	All	0.47	0/3274	0.71	1/4424 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	17	LYS	N-CA-C	6.08	127.42	111.00

There are no chirality outliers.

There are no planarity outliers.

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	A	1604	0	1577	240	0
1	B	1604	0	1579	194	1
2	A	25	0	28	4	0
2	B	25	0	26	9	0
3	A	4	0	5	3	0
4	A	23	0	0	5	0
4	B	17	0	0	0	0
All	All	3302	0	3215	414	1

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 64.

All (414) 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:A:108:CYS:SG	3:A:216:BME:S2	2.32	1.26
1:A:70:ARG:HH11	1:A:70:ARG:HG2	1.19	1.07
1:A:79:GLY:HA2	1:A:104:ARG:NH1	1.70	1.06
1:B:190:VAL:HG23	1:B:191:GLY:H	1.26	1.00
1:B:130:VAL:HG11	1:B:198:ILE:HG13	1.41	1.00
1:A:195:LEU:O	1:A:198:ILE:HG23	1.61	0.99
1:A:154:LYS:HG3	1:A:154:LYS:O	1.62	0.97
1:A:154:LYS:H	1:A:200:ARG:NH1	1.63	0.95
1:B:173:ASN:ND2	1:B:174:SER:H	1.66	0.94
1:A:195:LEU:H	1:A:198:ILE:HD13	1.32	0.94
1:A:69:ASP:O	1:A:70:ARG:HD3	1.69	0.93
1:A:52:GLN:NE2	1:A:82:ASP:H	1.66	0.93
1:A:52:GLN:HE22	1:A:82:ASP:N	1.67	0.92
1:A:70:ARG:HH11	1:A:70:ARG:CG	1.82	0.92
1:A:52:GLN:HE22	1:A:82:ASP:H	1.04	0.92
1:A:104:ARG:HG3	1:A:105:ILE:N	1.83	0.91
1:B:49:SER:OG	1:B:87:TYR:HB3	1.72	0.89
1:B:138:GLN:HG2	1:B:150:PRO:HD2	1.55	0.88
1:A:140:GLN:HE22	1:B:137:LYS:HA	1.38	0.88
1:A:59:LYS:HD3	1:A:70:ARG:HG3	1.55	0.87
1:A:195:LEU:N	1:A:198:ILE:HD13	1.90	0.86
1:B:173:ASN:HD22	1:B:174:SER:N	1.74	0.85
1:A:141:HIS:O	1:A:141:HIS:ND1	2.12	0.82
1:A:39:GLN:HG3	1:A:45:LEU:HG	1.61	0.82
1:B:138:GLN:HG3	1:B:147:PHE:CZ	2.14	0.81
1:A:118:ALA:HB1	4:A:224:HOH:O	1.79	0.81
1:A:108:CYS:SG	3:A:216:BME:C2	2.68	0.81
1:B:148:THR:O	1:B:149:GLY:O	2.00	0.80
1:A:138:GLN:HG3	1:A:147:PHE:HE1	1.47	0.80
1:B:82:ASP:OD2	1:B:99:ALA:HB1	1.82	0.80
1:A:138:GLN:HG2	1:B:21:LYS:HZ1	1.47	0.80
1:A:16:LEU:HD13	1:A:16:LEU:O	1.81	0.79
1:A:26:TRP:O	1:A:30:THR:HG22	1.82	0.79
1:B:113:GLN:HE21	1:B:113:GLN:HA	1.47	0.79
1:A:23:GLY:HA2	2:A:214:E64:O2	1.82	0.79
1:A:195:LEU:HD12	1:A:196:CYS:SG	2.23	0.79
1:A:79:GLY:HA2	1:A:104:ARG:HH11	1.48	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:VAL:O	1:A:35:GLU:HG3	1.83	0.78
1:B:155:LEU:HG	1:B:200:ARG:HH21	1.47	0.78
1:B:155:LEU:HD23	1:B:155:LEU:H	1.48	0.78
1:A:152:GLY:O	1:A:153:THR:O	2.02	0.78
1:A:79:GLY:HA2	1:A:104:ARG:HH12	1.49	0.77
1:B:41:ARG:HG3	1:B:207:LYS:HG2	1.66	0.76
1:B:138:GLN:HG3	1:B:147:PHE:CE1	2.21	0.76
1:A:169:TRP:NE1	1:A:189:ARG:HG2	2.01	0.75
1:A:154:LYS:H	1:A:200:ARG:HH12	1.32	0.75
1:B:52:GLN:HE22	1:B:82:ASP:H	1.31	0.74
1:A:140:GLN:NE2	1:B:140:GLN:OE1	2.20	0.74
1:A:14:ILE:HD12	1:A:47:SER:HB2	1.70	0.74
1:A:117:ASN:O	1:A:120:LYS:HD2	1.87	0.74
1:B:14:ILE:HG23	1:B:15:PRO:HD2	1.70	0.73
1:B:190:VAL:HG23	1:B:191:GLY:N	2.00	0.73
2:B:215:E64:H121	2:B:215:E64:N1	2.03	0.73
1:A:59:LYS:HD3	1:A:70:ARG:CG	2.18	0.73
1:A:138:GLN:HG3	1:A:147:PHE:CE1	2.24	0.73
1:A:16:LEU:HD21	1:A:177:ARG:HG2	1.71	0.73
1:A:116:GLU:OE2	1:A:198:ILE:HD11	1.89	0.73
1:B:31:VAL:O	1:B:35:GLU:HG3	1.88	0.73
1:B:131:ALA:HB1	1:B:155:LEU:HB3	1.68	0.72
2:A:214:E64:O1	1:B:137:LYS:NZ	2.23	0.72
1:A:71:ALA:O	1:A:75:ILE:HG13	1.90	0.72
1:B:4:HIS:HD2	1:B:164:TYR:O	1.72	0.72
1:A:70:ARG:HG2	1:A:70:ARG:NH1	1.97	0.72
1:A:81:ILE:HG13	1:A:82:ASP:N	2.04	0.72
1:A:79:GLY:CA	1:A:104:ARG:NH1	2.51	0.71
1:A:39:GLN:O	1:A:40:ILE:HG13	1.92	0.70
1:B:131:ALA:HB3	1:B:155:LEU:HD13	1.72	0.70
1:B:52:GLN:NE2	1:B:82:ASP:H	1.89	0.70
1:A:48:LEU:HD13	1:A:81:ILE:HG21	1.73	0.69
1:A:76:ILE:HD13	1:A:107:GLY:HA2	1.74	0.69
1:B:207:LYS:CG	1:B:207:LYS:O	2.39	0.69
1:A:79:GLY:CA	1:A:104:ARG:HH11	2.06	0.69
1:A:97:ARG:NH2	1:A:97:ARG:HG3	2.07	0.69
1:B:59:LYS:HG2	1:B:70:ARG:HG2	1.73	0.69
1:A:138:GLN:HG2	1:B:21:LYS:NZ	2.07	0.68
1:B:56:CYS:HB2	1:B:74:TYR:OH	1.92	0.68
1:B:65:GLY:HA3	2:B:215:E64:H132	1.76	0.68
1:A:97:ARG:HG3	1:A:97:ARG:HH21	1.56	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:LEU:HD11	1:B:200:ARG:HE	1.59	0.68
2:B:215:E64:H121	2:B:215:E64:C4	2.24	0.68
1:A:16:LEU:HD11	1:A:177:ARG:HE	1.59	0.67
1:A:48:LEU:HA	1:A:83:THR:HA	1.75	0.67
1:B:155:LEU:HD23	1:B:155:LEU:N	2.07	0.67
1:A:16:LEU:C	1:A:16:LEU:HD22	2.14	0.67
1:B:138:GLN:CG	1:B:150:PRO:HD2	2.25	0.67
1:B:173:ASN:HD22	1:B:174:SER:H	1.28	0.67
1:A:55:ASP:O	1:A:96:CYS:HB2	1.95	0.66
1:A:143:LYS:NZ	1:B:178:HIS:HB2	2.11	0.66
1:B:155:LEU:HG	1:B:200:ARG:NH2	2.10	0.66
1:A:76:ILE:CD1	1:A:107:GLY:HA2	2.26	0.66
1:A:13:VAL:CG1	1:A:172:ARG:NH1	2.59	0.65
1:A:108:CYS:SG	3:A:216:BME:H22	2.35	0.65
1:A:103:VAL:CG1	1:A:104:ARG:N	2.59	0.65
1:B:118:ALA:O	1:B:121:ASN:HB2	1.97	0.65
1:B:13:VAL:HG12	1:B:172:ARG:NH1	2.11	0.65
1:A:32:THR:HG22	4:A:221:HOH:O	1.97	0.64
1:B:150:PRO:O	1:B:151:CYS:HB3	1.96	0.64
1:A:159:VAL:HB	1:A:172:ARG:O	1.97	0.64
1:B:83:THR:HG23	1:B:102:VAL:HG21	1.79	0.64
1:A:147:PHE:CE1	1:A:149:GLY:HA3	2.33	0.64
1:B:16:LEU:O	1:B:177:ARG:NH1	2.31	0.64
1:A:16:LEU:CD2	1:A:177:ARG:HG2	2.28	0.63
2:B:215:E64:H121	2:B:215:E64:C3	2.27	0.63
1:A:103:VAL:HG12	1:A:104:ARG:N	2.13	0.63
1:A:195:LEU:H	1:A:198:ILE:CD1	2.09	0.63
1:B:207:LYS:HG2	1:B:207:LYS:O	1.98	0.63
1:A:104:ARG:CG	1:A:105:ILE:N	2.61	0.63
1:A:5:VAL:HG13	1:A:164:TYR:CZ	2.33	0.63
1:A:131:ALA:HB3	1:A:155:LEU:HD13	1.80	0.63
1:A:123:VAL:HG21	1:A:161:ILE:HG21	1.80	0.63
1:A:150:PRO:HG3	1:B:21:LYS:HG3	1.80	0.63
1:B:25:CYS:HB3	2:B:215:E64:O2	1.99	0.62
1:B:38:ASN:ND2	1:B:103:VAL:HG13	2.14	0.62
1:A:208:ALA:O	1:A:209:UNK:CB	2.47	0.62
1:A:143:LYS:NZ	1:B:178:HIS:CB	2.62	0.62
1:B:153:THR:HG21	1:B:193:CYS:SG	2.40	0.61
1:A:154:LYS:N	1:A:200:ARG:HH12	1.98	0.61
1:A:97:ARG:CG	1:A:98:ALA:H	2.12	0.61
1:A:116:GLU:OE2	1:A:194:GLY:HA2	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:PHE:CZ	2:B:215:E64:H91	2.35	0.61
1:A:140:GLN:NE2	1:B:137:LYS:HA	2.12	0.60
1:B:37:ILE:O	1:B:41:ARG:HG2	2.01	0.60
1:B:173:ASN:ND2	1:B:174:SER:N	2.36	0.60
1:A:16:LEU:CG	1:A:177:ARG:HG2	2.30	0.60
1:A:112:PRO:HA	4:A:220:HOH:O	2.01	0.60
1:A:148:THR:O	1:A:149:GLY:O	2.20	0.60
1:A:26:TRP:CG	1:A:66:GLY:HA3	2.36	0.60
1:A:97:ARG:HG2	1:A:98:ALA:H	1.65	0.60
1:B:51:GLN:HB2	1:B:89:TYR:HA	1.84	0.60
1:B:105:ILE:HB	1:B:206:THR:O	2.02	0.60
1:B:130:VAL:HG21	1:B:198:ILE:O	2.01	0.60
1:B:113:GLN:HE21	1:B:113:GLN:CA	2.13	0.60
1:A:178:HIS:HB2	1:B:143:LYS:HE3	1.83	0.60
1:B:177:ARG:O	1:B:183:GLY:N	2.34	0.60
1:A:33:THR:CG2	1:A:129:VAL:HG13	2.32	0.60
1:B:38:ASN:O	1:B:42:THR:OG1	2.18	0.59
1:A:97:ARG:HH21	1:A:97:ARG:CG	2.15	0.59
1:B:13:VAL:CG1	1:B:172:ARG:NH1	2.65	0.59
1:B:39:GLN:HB2	1:B:45:LEU:HD12	1.84	0.59
1:B:49:SER:O	1:B:81:ILE:HD11	2.02	0.59
1:A:143:LYS:HB2	4:A:222:HOH:O	2.03	0.59
1:B:14:ILE:HG22	1:B:15:PRO:O	2.02	0.58
1:A:141:HIS:NE2	1:B:179:TRP:HB2	2.17	0.58
1:A:1:LEU:HD11	1:A:121:ASN:OD1	2.03	0.58
1:B:84:GLU:OE2	1:B:88:PRO:HA	2.04	0.58
1:B:165:GLY:HA3	1:B:168:TYR:CE1	2.39	0.58
1:A:140:GLN:HE21	1:B:140:GLN:HB3	1.69	0.58
1:A:14:ILE:CD1	1:A:47:SER:HB2	2.33	0.57
1:A:123:VAL:HG21	1:A:161:ILE:CG2	2.34	0.57
1:A:167:ASP:O	1:A:188:LYS:HG3	2.04	0.57
1:B:7:TRP:HB2	1:B:162:VAL:HG13	1.86	0.57
1:A:19:GLN:HG2	1:A:174:SER:O	2.03	0.57
1:B:88:PRO:HG2	1:B:90:LYS:HE2	1.86	0.57
1:A:70:ARG:CG	1:A:70:ARG:NH1	2.52	0.57
1:A:168:TYR:CD1	1:A:168:TYR:C	2.77	0.57
1:B:182:GLN:O	1:B:182:GLN:HG2	2.04	0.56
1:A:14:ILE:HG22	1:A:15:PRO:O	2.04	0.56
1:A:116:GLU:HG2	1:A:199:ALA:HB2	1.88	0.56
1:A:169:TRP:CE2	1:A:189:ARG:HG2	2.39	0.56
1:B:59:LYS:NZ	1:B:70:ARG:HG2	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:TYR:CD1	1:A:164:TYR:C	2.79	0.56
1:B:4:HIS:CD2	1:B:164:TYR:O	2.56	0.56
1:B:56:CYS:O	1:B:58:LYS:HG3	2.06	0.56
1:B:138:GLN:HG3	1:B:147:PHE:HZ	1.66	0.55
1:A:81:ILE:HG13	1:A:82:ASP:H	1.71	0.55
1:A:1:LEU:HD12	1:A:120:LYS:HG3	1.89	0.55
1:B:72:TYR:HE1	1:B:205:PRO:HG3	1.72	0.55
1:B:8:ARG:HG3	1:B:162:VAL:HG11	1.88	0.55
1:A:32:THR:HA	1:A:35:GLU:OE1	2.07	0.55
1:A:132:ILE:O	1:A:155:LEU:HA	2.07	0.55
1:A:133:ASP:OD2	1:A:152:GLY:O	2.25	0.55
1:B:52:GLN:HE22	1:B:82:ASP:N	2.03	0.55
1:A:17:LYS:HB3	1:A:28:PHE:HE2	1.72	0.55
1:A:70:ARG:HH11	1:A:70:ARG:CB	2.20	0.55
1:B:38:ASN:HD22	1:B:103:VAL:HG13	1.70	0.55
1:A:116:GLU:OE1	1:A:190:VAL:N	2.38	0.54
1:B:54:VAL:HG13	1:B:62:GLY:CA	2.37	0.54
1:B:79:GLY:O	1:B:101:LYS:HG2	2.06	0.54
1:B:39:GLN:O	1:B:43:GLY:N	2.39	0.54
1:A:54:VAL:CG1	1:A:93:GLN:HA	2.38	0.54
1:A:132:ILE:O	1:A:156:ASN:N	2.39	0.54
1:B:159:VAL:HG21	1:B:171:VAL:HG11	1.90	0.54
1:A:167:ASP:N	1:A:167:ASP:OD1	2.39	0.53
1:A:17:LYS:NZ	1:A:84:GLU:OE1	2.41	0.53
1:A:33:THR:HG21	1:A:129:VAL:HG22	1.89	0.53
1:A:113:GLN:NE2	1:A:200:ARG:O	2.41	0.53
1:B:33:THR:HG21	1:B:129:VAL:CG1	2.39	0.53
1:A:26:TRP:O	1:A:30:THR:CG2	2.55	0.53
1:B:82:ASP:HB2	1:B:99:ALA:HB1	1.89	0.53
1:B:145:GLY:C	1:B:186:ARG:HH21	2.12	0.53
1:A:21:LYS:HE2	1:B:138:GLN:HB2	1.89	0.53
1:A:29:SER:HB2	1:A:129:VAL:HG12	1.91	0.53
1:A:55:ASP:OD2	1:A:95:PRO:O	2.26	0.53
1:A:157:HIS:NE2	1:A:173:ASN:ND2	2.56	0.53
1:B:145:GLY:CA	1:B:186:ARG:HH21	2.22	0.53
1:A:14:ILE:CG2	1:A:15:PRO:N	2.70	0.53
1:A:16:LEU:O	1:A:16:LEU:HD22	2.09	0.53
1:B:83:THR:CG2	1:B:102:VAL:HG11	2.38	0.53
1:A:138:GLN:HB2	1:B:21:LYS:HE2	1.91	0.52
1:B:132:ILE:O	1:B:155:LEU:HA	2.08	0.52
1:A:38:ASN:O	1:A:42:THR:OG1	2.17	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ASP:HB3	1:A:95:PRO:O	2.09	0.52
1:B:15:PRO:HB2	1:B:177:ARG:NH2	2.25	0.52
1:A:123:VAL:O	1:A:123:VAL:CG1	2.57	0.52
1:A:148:THR:CG2	1:A:149:GLY:N	2.72	0.52
1:B:165:GLY:O	1:B:167:ASP:N	2.42	0.52
1:A:141:HIS:HD1	1:A:141:HIS:C	2.13	0.52
1:A:49:SER:CB	1:A:87:TYR:HB3	2.39	0.52
1:A:139:PHE:O	1:A:179:TRP:HZ2	1.92	0.52
1:B:25:CYS:HB2	1:B:157:HIS:CE1	2.44	0.52
1:B:49:SER:OG	1:B:87:TYR:O	2.28	0.52
1:A:138:GLN:HB2	1:B:21:LYS:CE	2.40	0.52
1:A:28:PHE:CD1	1:A:50:GLU:HG2	2.46	0.51
1:A:49:SER:O	1:A:81:ILE:HD11	2.11	0.51
1:A:126:GLN:NE2	1:A:204:TYR:CD1	2.78	0.51
1:A:140:GLN:HE21	1:B:140:GLN:CB	2.23	0.51
1:A:14:ILE:HG23	1:A:15:PRO:CD	2.40	0.51
1:A:198:ILE:HG12	1:A:199:ALA:N	2.25	0.51
1:A:118:ALA:O	1:A:121:ASN:HB2	2.10	0.51
1:A:23:GLY:CA	2:A:214:E64:O2	2.55	0.51
1:A:25:CYS:SG	1:A:26:TRP:N	2.84	0.51
1:A:54:VAL:O	1:A:93:GLN:HG3	2.11	0.51
1:A:97:ARG:CG	1:A:98:ALA:N	2.74	0.51
1:A:150:PRO:HD2	1:B:21:LYS:HZ1	1.76	0.51
1:A:76:ILE:HD11	1:A:107:GLY:O	2.11	0.51
1:A:97:ARG:NH2	1:A:97:ARG:CG	2.74	0.50
1:A:153:THR:HB	1:A:200:ARG:HH11	1.76	0.50
1:A:13:VAL:HG13	1:A:172:ARG:NH1	2.25	0.50
1:A:138:GLN:CG	1:A:147:PHE:HE1	2.20	0.50
1:B:59:LYS:HZ3	1:B:70:ARG:HG2	1.74	0.50
1:B:15:PRO:O	1:B:172:ARG:NH2	2.40	0.50
1:A:142:TYR:CD2	1:A:147:PHE:HB2	2.46	0.50
1:A:169:TRP:N	1:A:187:MET:O	2.39	0.50
1:A:154:LYS:H	1:A:200:ARG:HH11	1.53	0.50
1:A:140:GLN:OE1	1:B:137:LYS:HG2	2.12	0.50
1:B:14:ILE:HG23	1:B:15:PRO:CD	2.41	0.50
1:A:188:LYS:HZ2	1:A:188:LYS:CB	2.25	0.50
1:B:16:LEU:HD21	1:B:184:TYR:CZ	2.47	0.50
1:B:68:PHE:CD2	1:B:203:PHE:HD2	2.30	0.50
1:A:40:ILE:HG22	1:A:41:ARG:HG2	1.92	0.50
1:A:32:THR:HA	1:A:35:GLU:CD	2.32	0.49
1:B:26:TRP:CD1	1:B:66:GLY:N	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:ASP:OD2	1:A:208:ALA:HB2	2.11	0.49
1:A:111:VAL:HG12	1:A:112:PRO:HD2	1.94	0.49
1:A:131:ALA:HA	1:A:157:HIS:O	2.13	0.49
1:B:36:SER:O	1:B:40:ILE:HG13	2.12	0.49
1:A:21:LYS:HG3	1:B:150:PRO:HG3	1.93	0.49
1:A:83:THR:HG23	1:A:102:VAL:HG21	1.94	0.49
1:B:16:LEU:CD1	1:B:183:GLY:HA3	2.43	0.49
1:A:150:PRO:O	1:A:151:CYS:HB3	2.12	0.49
1:A:54:VAL:HG22	1:A:62:GLY:HA2	1.94	0.49
1:A:195:LEU:O	1:A:198:ILE:CG2	2.49	0.49
1:B:145:GLY:HA2	1:B:186:ARG:HH21	1.78	0.49
1:A:116:GLU:OE1	1:A:190:VAL:HA	2.12	0.49
1:B:89:TYR:CZ	1:B:91:ALA:HA	2.48	0.49
1:A:41:ARG:HD2	1:A:207:LYS:HB3	1.95	0.48
1:A:105:ILE:HG22	1:A:206:THR:O	2.13	0.48
1:B:159:VAL:HG21	1:B:171:VAL:CG1	2.43	0.48
1:B:179:TRP:HH2	1:B:185:THR:HG21	1.77	0.48
1:A:33:THR:HG21	1:A:129:VAL:HG13	1.94	0.48
1:A:178:HIS:ND1	1:A:178:HIS:N	2.58	0.48
1:B:14:ILE:HD12	1:B:47:SER:HB2	1.94	0.48
1:B:123:VAL:HG21	1:B:161:ILE:HG21	1.95	0.48
1:A:14:ILE:HG23	1:A:15:PRO:HD2	1.95	0.48
1:A:16:LEU:CD1	1:A:16:LEU:N	2.77	0.48
1:B:33:THR:HG21	1:B:129:VAL:HG13	1.96	0.48
1:B:82:ASP:OD1	1:B:86:ASN:HB2	2.12	0.48
1:A:141:HIS:O	1:A:141:HIS:CG	2.65	0.48
1:A:78:ASN:CG	1:A:78:ASN:O	2.52	0.48
1:A:20:GLY:C	1:A:22:CYS:H	2.16	0.48
1:A:89:TYR:CE1	1:A:91:ALA:HA	2.49	0.48
1:B:105:ILE:HD13	1:B:105:ILE:O	2.14	0.48
1:B:145:GLY:HA2	1:B:186:ARG:NH2	2.29	0.48
1:A:154:LYS:O	1:A:154:LYS:CG	2.48	0.48
2:B:215:E64:C12	2:B:215:E64:HN3	2.24	0.48
1:A:97:ARG:CD	1:A:98:ALA:H	2.27	0.47
1:A:165:GLY:O	1:A:189:ARG:NH2	2.47	0.47
1:B:25:CYS:CB	1:B:157:HIS:CE1	2.96	0.47
1:A:103:VAL:CG1	1:A:104:ARG:H	2.27	0.47
1:B:78:ASN:O	1:B:101:LYS:HE2	2.14	0.47
1:A:46:ILE:HD12	1:A:47:SER:H	1.79	0.47
1:A:20:GLY:HA2	4:A:226:HOH:O	2.15	0.47
1:A:123:VAL:O	1:A:123:VAL:HG12	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:CYS:SG	1:A:195:LEU:CD1	3.03	0.47
1:B:30:THR:O	1:B:34:VAL:HG23	2.14	0.47
1:A:97:ARG:CD	1:A:98:ALA:N	2.78	0.47
1:A:153:THR:O	1:A:154:LYS:CB	2.62	0.47
1:B:136:SER:O	1:B:140:GLN:HB2	2.14	0.47
1:B:151:CYS:SG	1:B:151:CYS:O	2.72	0.47
1:A:70:ARG:NH1	1:A:70:ARG:CB	2.78	0.47
1:B:31:VAL:HG11	1:B:50:GLU:OE2	2.15	0.46
1:B:68:PHE:HD2	1:B:203:PHE:CD2	2.33	0.46
1:B:111:VAL:HB	1:B:202:PRO:HG2	1.97	0.46
1:B:177:ARG:HA	1:B:182:GLN:C	2.36	0.46
1:B:177:ARG:O	1:B:178:HIS:C	2.54	0.46
1:B:48:LEU:HD13	1:B:81:ILE:HG21	1.96	0.46
1:A:21:LYS:HD2	1:A:21:LYS:N	2.30	0.46
1:A:175:TRP:CE3	1:B:141:HIS:NE2	2.82	0.46
1:B:56:CYS:CB	1:B:74:TYR:OH	2.60	0.46
1:A:156:ASN:C	2:A:214:E64:H72	2.36	0.46
1:A:16:LEU:HG	1:A:177:ARG:HG2	1.95	0.46
1:B:65:GLY:HA2	2:B:215:E64:O3	2.15	0.45
1:A:5:VAL:CG1	1:A:164:TYR:CZ	2.98	0.45
1:A:69:ASP:OD1	1:A:203:PHE:CE2	2.69	0.45
1:B:113:GLN:NE2	1:B:200:ARG:O	2.49	0.45
1:A:105:ILE:CG2	1:A:206:THR:O	2.65	0.45
1:B:16:LEU:HD12	1:B:183:GLY:HA3	1.98	0.45
1:A:4:HIS:CD2	1:A:164:TYR:O	2.70	0.45
1:A:54:VAL:HG12	1:A:93:GLN:HA	1.98	0.45
1:B:107:GLY:C	1:B:206:THR:CG2	2.85	0.45
1:B:201:LEU:HA	1:B:201:LEU:HD23	1.71	0.45
1:B:48:LEU:HD22	1:B:81:ILE:HG23	1.97	0.45
1:B:155:LEU:CG	1:B:200:ARG:HH21	2.25	0.45
1:A:16:LEU:HD11	1:A:177:ARG:NE	2.30	0.45
1:A:116:GLU:HA	1:A:119:LEU:HB3	1.98	0.45
1:A:136:SER:OG	1:A:139:PHE:CB	2.64	0.45
1:B:130:VAL:HG13	1:B:131:ALA:N	2.31	0.45
1:A:39:GLN:HG3	1:A:45:LEU:CG	2.41	0.45
1:A:14:ILE:HD12	1:A:47:SER:CB	2.44	0.44
1:A:116:GLU:OE1	1:A:190:VAL:CA	2.65	0.44
1:B:130:VAL:CG2	1:B:198:ILE:O	2.65	0.44
1:B:204:TYR:HA	1:B:205:PRO:HD3	1.78	0.44
1:A:16:LEU:HD23	1:A:176:GLY:C	2.36	0.44
1:A:129:VAL:N	1:A:203:PHE:O	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:GLN:NE2	1:B:140:GLN:CB	2.80	0.44
1:B:21:LYS:HD2	1:B:21:LYS:N	2.32	0.44
1:A:38:ASN:ND2	1:A:103:VAL:HG22	2.32	0.44
1:A:107:GLY:O	1:A:108:CYS:HB3	2.17	0.44
1:A:111:VAL:HB	1:A:202:PRO:CG	2.47	0.44
1:A:19:GLN:HA	1:A:28:PHE:CZ	2.53	0.44
1:B:1:LEU:HA	1:B:2:PRO:HD3	1.75	0.44
1:A:159:VAL:HG21	1:A:171:VAL:CG1	2.48	0.43
1:B:39:GLN:HG3	1:B:45:LEU:HD13	2.00	0.43
1:B:190:VAL:CG2	1:B:191:GLY:N	2.70	0.43
1:A:3:GLU:O	1:A:164:TYR:HE1	2.00	0.43
1:A:78:ASN:O	1:A:101:LYS:HE2	2.18	0.43
1:B:54:VAL:HG13	1:B:62:GLY:HA2	1.99	0.43
1:B:105:ILE:N	1:B:105:ILE:CD1	2.81	0.43
1:A:164:TYR:C	1:A:164:TYR:HD1	2.21	0.43
1:B:147:PHE:CG	1:B:148:THR:N	2.80	0.43
1:B:8:ARG:CG	1:B:162:VAL:HG11	2.48	0.43
1:B:155:LEU:H	1:B:155:LEU:CD2	2.07	0.43
1:B:166:LYS:H	1:B:166:LYS:HG3	1.50	0.43
1:A:20:GLY:O	1:A:22:CYS:N	2.44	0.43
1:A:37:ILE:HD12	1:A:37:ILE:HA	1.86	0.43
1:A:49:SER:OG	1:A:87:TYR:HB3	2.18	0.43
1:B:138:GLN:HB3	1:B:150:PRO:HG2	2.00	0.43
1:A:14:ILE:HG23	1:A:15:PRO:N	2.33	0.43
1:A:169:TRP:O	1:A:186:ARG:HA	2.18	0.43
1:A:33:THR:HG23	1:A:129:VAL:HG13	2.00	0.43
1:B:148:THR:O	1:B:149:GLY:C	2.56	0.43
1:B:155:LEU:HD11	1:B:200:ARG:HB3	2.01	0.43
1:B:54:VAL:O	1:B:93:GLN:NE2	2.43	0.43
1:A:72:TYR:OH	1:A:205:PRO:HD3	2.19	0.43
1:B:82:ASP:CB	1:B:99:ALA:HB1	2.49	0.43
1:B:177:ARG:HG3	1:B:177:ARG:HH11	1.84	0.43
1:B:83:THR:HG22	1:B:102:VAL:HG11	2.00	0.42
1:B:131:ALA:CB	1:B:155:LEU:HB3	2.44	0.42
1:B:162:VAL:HG12	1:B:162:VAL:O	2.17	0.42
1:B:115:ASN:OD1	1:B:118:ALA:CB	2.67	0.42
1:B:131:ALA:CB	2:B:215:E64:C9	2.98	0.42
1:B:72:TYR:CE2	1:B:203:PHE:HB3	2.54	0.42
1:B:104:ARG:O	1:B:207:LYS:NZ	2.41	0.42
1:A:1:LEU:CD1	1:A:120:LYS:HG3	2.48	0.42
1:B:28:PHE:CG	1:B:50:GLU:HG2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:GLN:O	1:B:43:GLY:HA2	2.19	0.42
1:A:13:VAL:CG1	1:A:172:ARG:HH11	2.33	0.42
1:A:126:GLN:NE2	1:A:205:PRO:O	2.52	0.42
1:B:5:VAL:O	1:B:163:GLY:HA3	2.20	0.42
1:B:105:ILE:HD13	1:B:105:ILE:H	1.84	0.42
1:B:8:ARG:HD3	1:B:184:TYR:CE2	2.55	0.42
1:A:32:THR:HA	1:A:35:GLU:CG	2.49	0.42
1:B:37:ILE:HD13	1:B:37:ILE:HA	1.81	0.42
1:A:64:LYS:HB3	1:A:64:LYS:HE3	1.83	0.42
1:A:143:LYS:HZ3	1:B:178:HIS:HB2	1.83	0.42
1:B:188:LYS:HG2	1:B:189:ARG:H	1.85	0.42
1:A:32:THR:HG21	1:A:172:ARG:NE	2.35	0.41
1:A:131:ALA:HB1	1:A:155:LEU:HB3	2.01	0.41
1:B:32:THR:HA	1:B:35:GLU:OE1	2.20	0.41
1:A:116:GLU:OE2	1:A:194:GLY:CA	2.68	0.41
1:A:137:LYS:HZ3	1:B:175:TRP:HE1	1.68	0.41
1:A:189:ARG:O	1:A:190:VAL:HB	2.20	0.41
1:B:31:VAL:CG1	1:B:50:GLU:OE2	2.68	0.41
1:B:33:THR:HG21	1:B:129:VAL:HG12	2.02	0.41
1:B:184:TYR:CD1	1:B:184:TYR:N	2.88	0.41
1:A:1:LEU:CD1	1:A:120:LYS:CG	2.99	0.41
1:A:157:HIS:HD2	1:A:159:VAL:HG13	1.85	0.41
1:B:68:PHE:HB2	1:B:203:PHE:CD2	2.55	0.41
1:B:81:ILE:HG13	1:B:82:ASP:N	2.36	0.41
1:B:148:THR:HA	1:B:195:LEU:CD2	2.50	0.41
1:B:150:PRO:O	1:B:151:CYS:CB	2.64	0.41
1:B:190:VAL:CG2	1:B:191:GLY:H	2.04	0.41
1:A:29:SER:HB2	1:A:129:VAL:CG1	2.50	0.41
1:A:140:GLN:NE2	1:B:140:GLN:HB2	2.35	0.41
1:A:169:TRP:CD1	1:A:189:ARG:HG2	2.55	0.41
1:A:82:ASP:OD2	1:A:99:ALA:HB1	2.21	0.41
1:A:61:HIS:O	1:A:62:GLY:C	2.59	0.41
1:A:120:LYS:O	1:A:124:ALA:HB2	2.21	0.41
1:B:29:SER:O	1:B:129:VAL:HG12	2.21	0.41
1:B:54:VAL:HG13	1:B:62:GLY:HA3	2.02	0.41
1:B:90:LYS:O	1:B:91:ALA:HB3	2.21	0.41
1:B:141:HIS:O	1:B:142:TYR:C	2.60	0.41
1:A:17:LYS:O	1:A:18:ASN:HB2	2.21	0.41
1:A:150:PRO:CD	1:B:21:LYS:HZ1	2.33	0.41
1:A:103:VAL:HG13	1:A:104:ARG:H	1.86	0.40
1:A:90:LYS:O	1:A:91:ALA:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:GLN:O	1:A:114:CYS:C	2.59	0.40
1:B:39:GLN:O	1:B:43:GLY:CA	2.69	0.40
1:B:134:ALA:HB1	1:B:139:PHE:CE1	2.55	0.40
1:B:177:ARG:O	1:B:179:TRP:O	2.39	0.40
1:B:53:LEU:HD23	1:B:74:TYR:CD2	2.57	0.40
1:B:169:TRP:O	1:B:186:ARG:HA	2.21	0.40
1:A:171:VAL:HB	1:A:185:THR:HG23	2.02	0.40
1:B:78:ASN:O	1:B:101:LYS:CE	2.70	0.40

All (1) 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)
1:B:13:VAL:O	1:B:70:ARG:NH2[1_655]	2.05	0.15

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	207/209 (99%)	163 (79%)	28 (14%)	16 (8%)	1 2
1	B	207/209 (99%)	164 (79%)	32 (16%)	11 (5%)	2 5
All	All	414/418 (99%)	327 (79%)	60 (14%)	27 (6%)	1 3

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	40	ILE
1	A	149	GLY
1	A	153	THR
1	B	97	ARG

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Mol	Chain	Res	Type
1	B	148	THR
1	B	149	GLY
1	B	151	CYS
1	B	166	LYS
1	B	190	VAL
1	A	62	GLY
1	A	114	CYS
1	A	151	CYS
1	A	190	VAL
1	B	41	ARG
1	B	194	GLY
1	B	208	ALA
1	A	70	ARG
1	A	108	CYS
1	A	21	LYS
1	A	166	LYS
1	A	167	ASP
1	A	198	ILE
1	B	202	PRO
1	B	198	ILE
1	A	145	GLY
1	A	191	GLY

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	166/166 (100%)	130 (78%)	36 (22%)	1 2
1	B	166/166 (100%)	135 (81%)	31 (19%)	1 3
All	All	332/332 (100%)	265 (80%)	67 (20%)	1 3

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ILE

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Mol	Chain	Res	Type
1	A	16	LEU
1	A	21	LYS
1	A	41	ARG
1	A	45	LEU
1	A	53	LEU
1	A	55	ASP
1	A	67	TYR
1	A	69	ASP
1	A	70	ARG
1	A	73	GLN
1	A	81	ILE
1	A	87	TYR
1	A	90	LYS
1	A	92	PHE
1	A	97	ARG
1	A	100	LYS
1	A	104	ARG
1	A	105	ILE
1	A	109	LYS
1	A	112	PRO
1	A	113	GLN
1	A	129	VAL
1	A	142	TYR
1	A	154	LYS
1	A	155	LEU
1	A	156	ASN
1	A	166	LYS
1	A	173	ASN
1	A	177	ARG
1	A	178	HIS
1	A	189	ARG
1	A	195	LEU
1	A	198	ILE
1	A	201	LEU
1	A	206	THR
1	B	3	GLU
1	B	5	VAL
1	B	16	LEU
1	B	21	LYS
1	B	36	SER
1	B	37	ILE
1	B	49	SER

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Mol	Chain	Res	Type
1	B	57	SER
1	B	59	LYS
1	B	67	TYR
1	B	87	TYR
1	B	92	PHE
1	B	97	ARG
1	B	105	ILE
1	B	106	ASP
1	B	112	PRO
1	B	113	GLN
1	B	125	SER
1	B	126	GLN
1	B	130	VAL
1	B	135	SER
1	B	136	SER
1	B	148	THR
1	B	154	LYS
1	B	155	LEU
1	B	166	LYS
1	B	173	ASN
1	B	185	THR
1	B	200	ARG
1	B	206	THR
1	B	207	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	38	ASN
1	A	39	GLN
1	A	52	GLN
1	A	86	ASN
1	A	113	GLN
1	A	126	GLN
1	A	140	GLN
1	A	173	ASN
1	B	4	HIS
1	B	38	ASN
1	B	52	GLN
1	B	113	GLN
1	B	173	ASN

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Mol	Chain	Res	Type
1	B	182	GLN

5.3.3 RNA

There are no RNA molecules 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 monosaccharides in this entry.

5.6 Ligand geometry

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	BME	A	216	-	3,3,3	0.28	0	1,2,2	0.52	0
2	E64	A	214	-	24,24,24	0.76	2 (8%)	27,30,30	1.58	6 (22%)
2	E64	B	215	-	24,24,24	0.76	2 (8%)	27,30,30	1.82	3 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BME	A	216	-	-	0/1/1/1	-
2	E64	B	215	-	1/1/7/10	15/29/29/29	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	E64	A	214	-	-	15/29/29/29	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	215	E64	O1-C1	2.65	1.30	1.22
2	A	214	E64	O1-C1	2.56	1.30	1.22
2	A	214	E64	O2-C1	-2.26	1.23	1.30
2	B	215	E64	O2-C1	-2.19	1.23	1.30

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	215	E64	O3-C3-C4	7.06	127.74	111.01
2	A	214	E64	O3-C3-C2	4.92	122.13	110.05
2	B	215	E64	O3-C3-C2	4.42	120.91	110.05
2	A	214	E64	N3-C16-N5	3.80	127.38	120.70
2	A	214	E64	N4-C16-N3	-2.65	113.08	119.19
2	A	214	E64	O3-C3-C4	2.52	116.98	111.01
2	A	214	E64	O2-C1-C2	2.16	120.99	114.07
2	B	215	E64	O2-C1-C2	2.15	120.95	114.07
2	A	214	E64	O1-C1-C2	-2.10	116.06	122.80

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	215	E64	C3

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	214	E64	O3-C3-C4-N1
2	A	214	E64	N4-C16-N3-C15
2	B	215	E64	C1-C2-C3-O3
2	B	215	E64	O3-C3-C4-O4
2	B	215	E64	C3-C4-N1-C6
2	B	215	E64	C6-C11-N2-C12
2	B	215	E64	O5-C11-N2-C12
2	B	215	E64	N4-C16-N3-C15
2	B	215	E64	N5-C16-N3-C15
2	B	215	E64	O4-C4-N1-C6

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Mol	Chain	Res	Type	Atoms
2	A	214	E64	C12-C13-C14-C15
2	A	214	E64	N1-C6-C7-C8
2	A	214	E64	N2-C12-C13-C14
2	A	214	E64	C11-C6-C7-C8
2	B	215	E64	N2-C12-C13-C14
2	A	214	E64	C6-C7-C8-C9
2	A	214	E64	C6-C7-C8-C10
2	B	215	E64	C13-C14-C15-N3
2	A	214	E64	C13-C14-C15-N3
2	A	214	E64	O3-C3-C4-O4
2	A	214	E64	N5-C16-N3-C15
2	B	215	E64	O5-C11-C6-C7
2	B	215	E64	N2-C11-C6-C7
2	B	215	E64	O3-C3-C4-N1
2	A	214	E64	C2-C3-C4-O4
2	B	215	E64	C2-C3-C4-O4
2	A	214	E64	O5-C11-C6-C7
2	A	214	E64	N2-C11-C6-C7
2	A	214	E64	C2-C3-C4-N1
2	B	215	E64	C2-C3-C4-N1

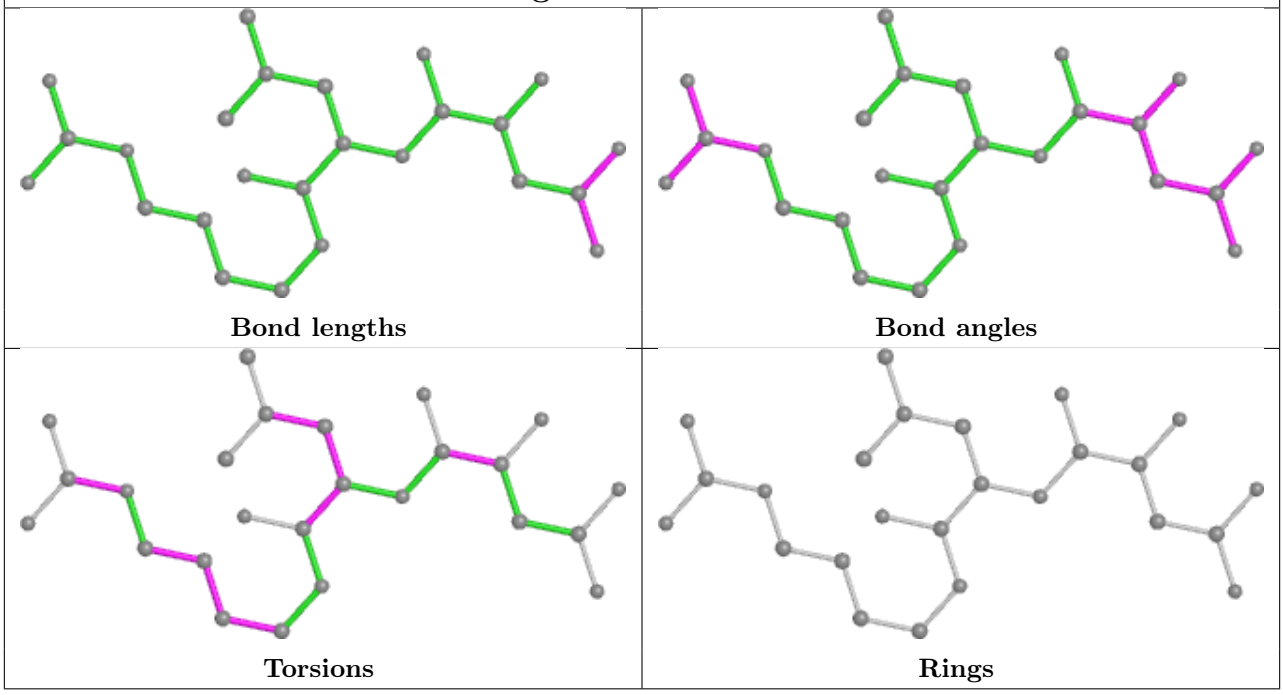
There are no ring outliers.

3 monomers are involved in 16 short contacts:

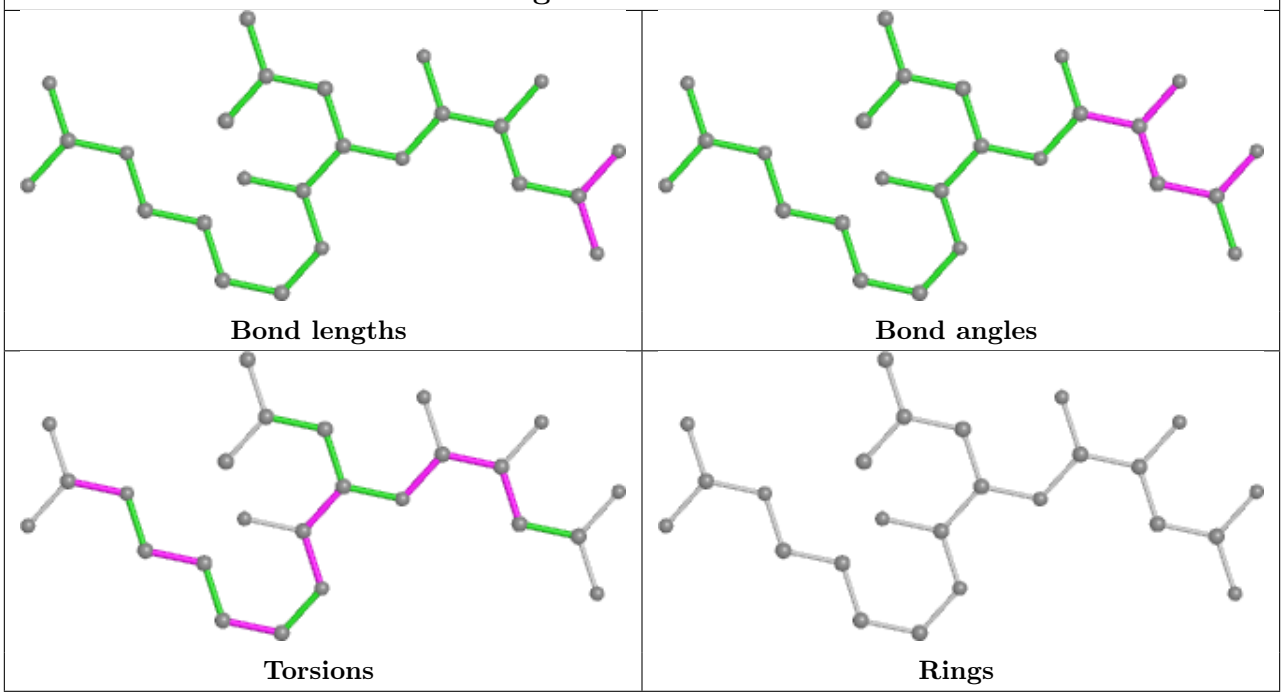
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	216	BME	3	0
2	A	214	E64	4	0
2	B	215	E64	9	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand E64 A 214



Ligand E64 B 215



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^{th} 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	208/209 (99%)	-0.20	0	100	100	12, 12, 12, 16	0
1	B	208/209 (99%)	-0.25	0	100	100	12, 12, 12, 16	0
All	All	416/418 (99%)	-0.23	0	100	100	12, 12, 12, 16	0

There are no RSRZ outliers to report.

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 monosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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.

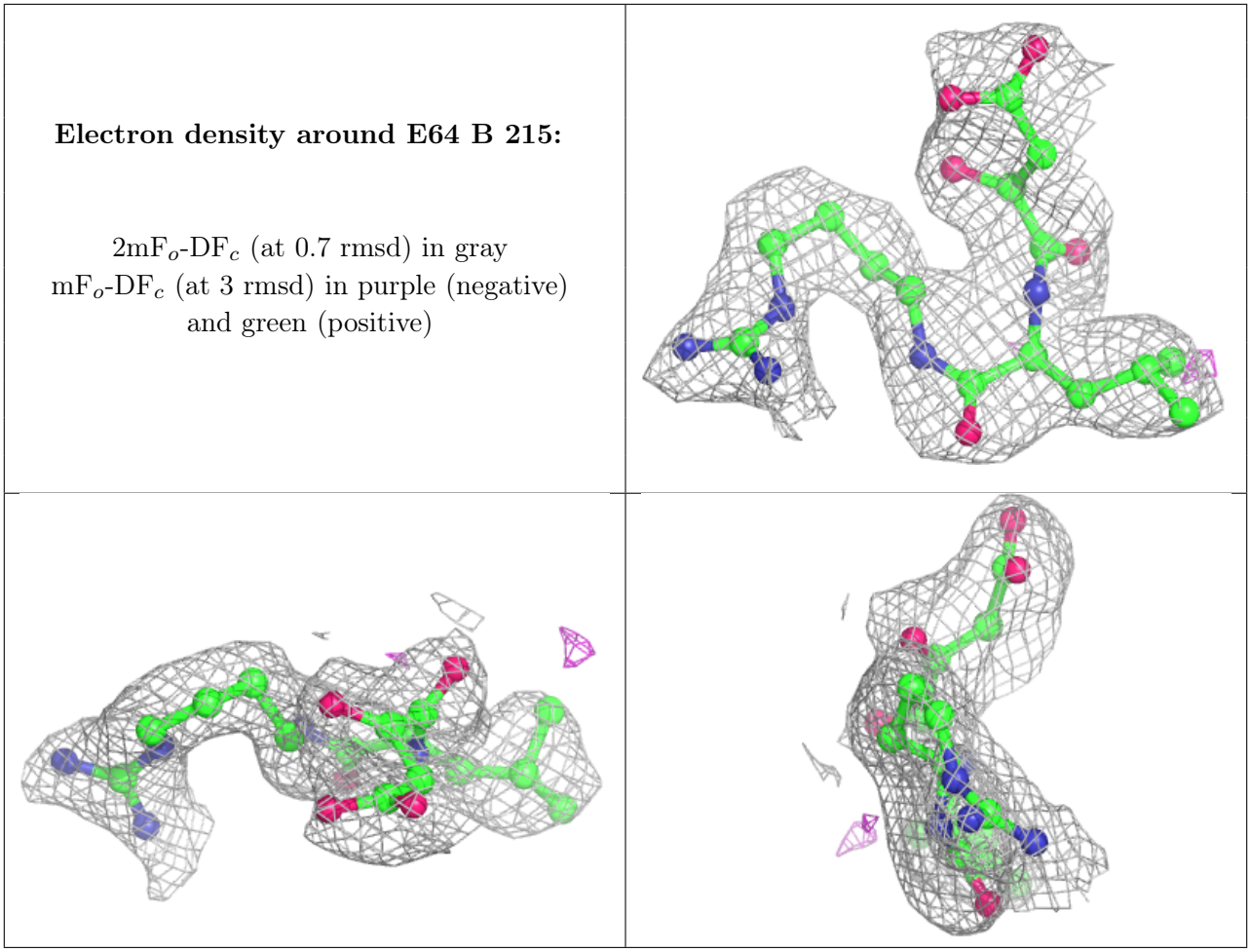
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	BME	A	216	4/4	0.85	0.22	45,50,51,52	0
2	E64	B	215	25/25	0.87	0.23	4,15,22,25	0
2	E64	A	214	25/25	0.89	0.20	1,12,19,22	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different

orientation to approximate a three-dimensional view.

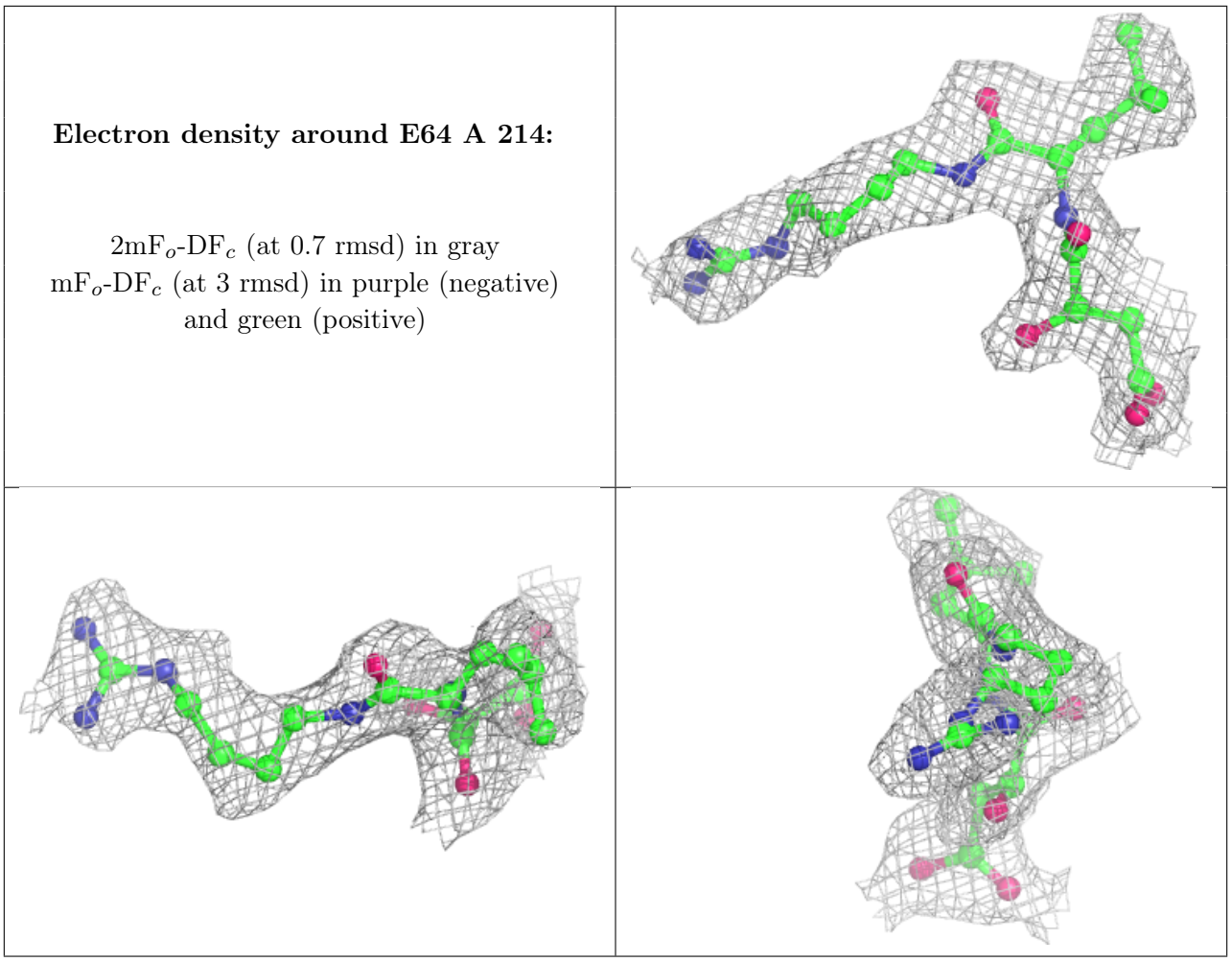
Electron density around E64 B 215:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around E64 A 214:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.