



# Full wwPDB X-ray Structure Validation Report i

May 26, 2020 – 08:36 am BST

PDB ID : 5OY6  
Title : Crystal structure of the ACVR1 (ALK2) kinase in complex with cyclical inhibitor OD36.  
Authors : Williams, E.P.; Pinkas, D.M.; Krojer, T.; Kupinska, K.; Mahajan, P.; Burgess-Brown, N.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Bullock, A.N.  
Deposited on : 2017-09-07  
Resolution : 2.56 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the i symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

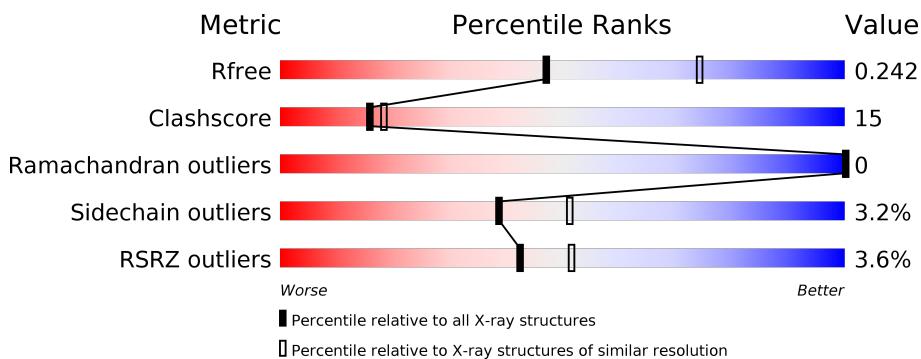
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## X-RAY DIFFRACTION

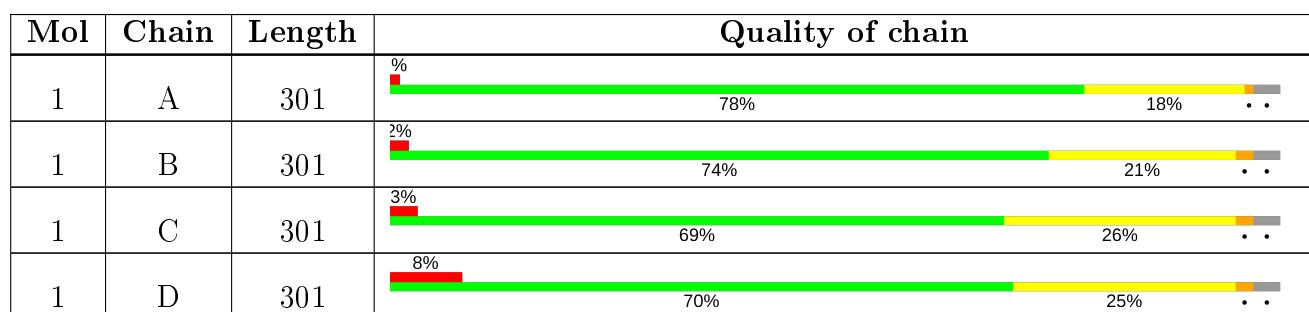
The reported resolution of this entry is 2.56 Å.

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	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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 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.



## 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 9055 atoms, of which 12 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.

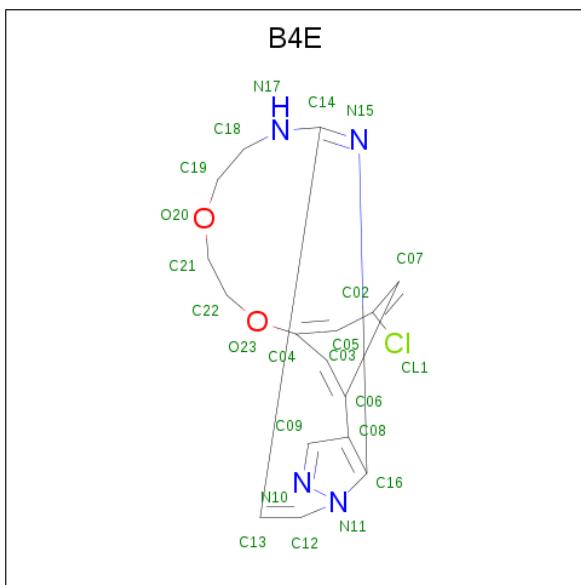
- Molecule 1 is a protein called Activin receptor type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total 2269	C 1445	N 389	O 420	S 15	0	0	0
1	B	292	Total 2263	C 1445	N 389	O 414	S 15	0	0	0
1	C	291	Total 2203	C 1410	N 373	O 405	S 15	0	0	0
1	D	292	Total 2138	C 1367	N 365	O 392	S 14	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

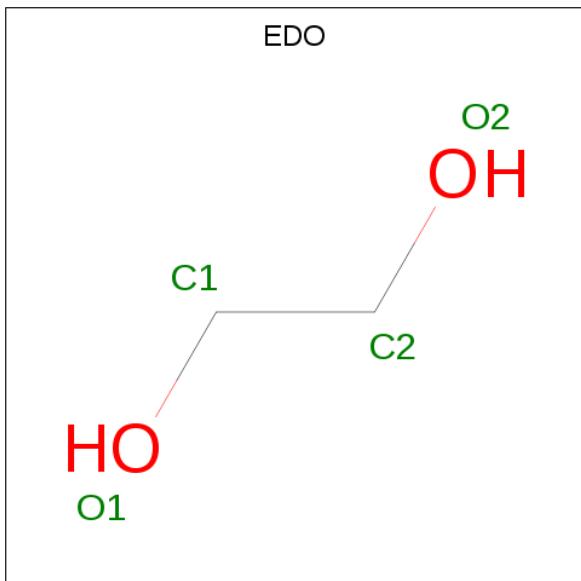
Chain	Residue	Modelled	Actual	Comment	Reference
A	199	SER	-	expression tag	UNP Q04771
A	200	MET	-	expression tag	UNP Q04771
A	207	ASP	GLN	engineered mutation	UNP Q04771
B	199	SER	-	expression tag	UNP Q04771
B	200	MET	-	expression tag	UNP Q04771
B	207	ASP	GLN	engineered mutation	UNP Q04771
C	199	SER	-	expression tag	UNP Q04771
C	200	MET	-	expression tag	UNP Q04771
C	207	ASP	GLN	engineered mutation	UNP Q04771
D	199	SER	-	expression tag	UNP Q04771
D	200	MET	-	expression tag	UNP Q04771
D	207	ASP	GLN	engineered mutation	UNP Q04771

- Molecule 2 is cyclical inhibitor OD36 (three-letter code: B4E) (formula: C<sub>16</sub>H<sub>15</sub>ClN<sub>4</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total		C	Cl	N	O	
			23		16	1	4	2	0
2	B	1	Total		C	Cl	N	O	
			23		16	1	4	2	0
2	C	1	Total		C	Cl	N	O	
			23		16	1	4	2	0
2	D	1	Total		C	Cl	N	O	
			23		16	1	4	2	0

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H O 10 2 6 2	0	0
3	A	1	Total C H O 10 2 6 2	0	0

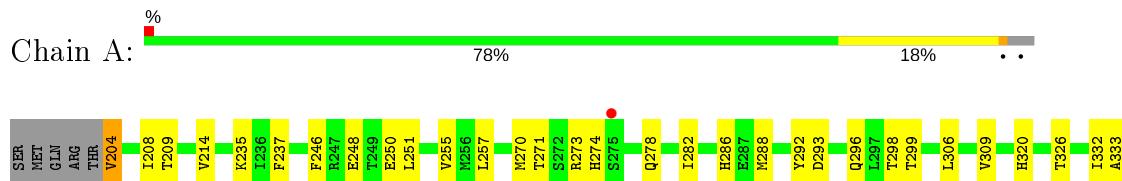
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	41	Total O 41 41	0	0
4	B	16	Total O 16 16	0	0
4	C	7	Total O 7 7	0	0
4	D	6	Total O 6 6	0	0

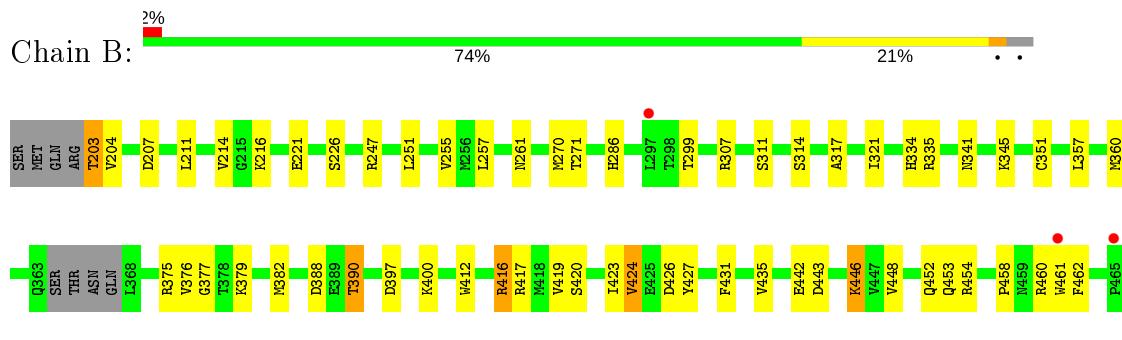
### 3 Residue-property plots

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

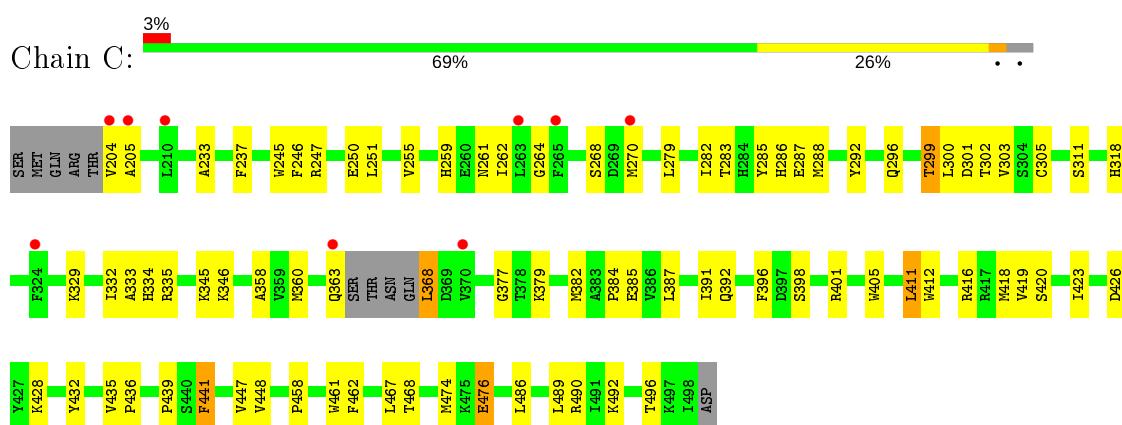
- Molecule 1: Activin receptor type-1



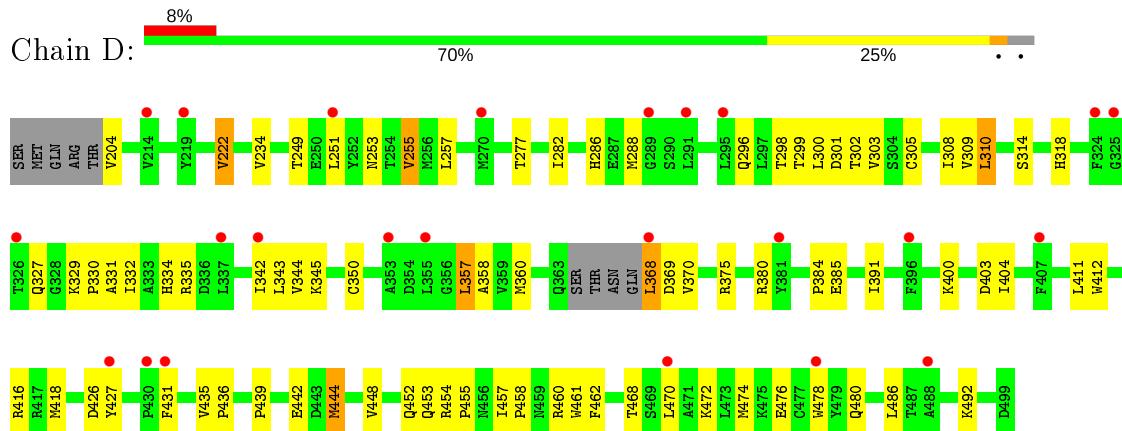
- Molecule 1: Activin receptor type-1



- Molecule 1: Activin receptor type-1



- Molecule 1: Activin receptor type-1



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.26 Å    103.53 Å    84.84 Å 90.00°    116.52°    90.00°	Depositor
Resolution (Å)	71.90 – 2.56 71.90 – 2.56	Depositor EDS
% Data completeness (in resolution range)	99.7 (71.90-2.56) 99.8 (71.90-2.56)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle^1$	2.32 (at 2.55 Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ????)	Depositor
$R$ , $R_{free}$	0.211 , 0.242 0.212 , 0.242	Depositor DCC
$R_{free}$ test set	2125 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	62.3	Xtriage
Anisotropy	0.770	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 61.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9055	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: B4E, EDO

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	0.40	0/2320	0.57	0/3158
1	B	0.35	0/2314	0.54	0/3148
1	C	0.33	0/2254	0.51	0/3078
1	D	0.32	0/2187	0.52	1/2992 (0.0%)
All	All	0.35	0/9075	0.53	1/12376 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	327	GLN	N-CA-C	-5.49	96.18	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts i

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2269	0	2185	51	0
1	B	2263	0	2190	48	0
1	C	2203	0	2058	84	0
1	D	2138	0	1948	71	0
2	A	23	0	0	2	0
2	B	23	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	23	0	0	0	0
2	D	23	0	0	1	0
3	A	8	12	12	0	0
4	A	41	0	0	8	0
4	B	16	0	0	2	0
4	C	7	0	0	2	0
4	D	6	0	0	3	0
All	All	9043	12	8393	254	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (254) 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:435:VAL:CG2	1:A:439:PRO:HB3	1.90	0.99
1:C:302:THR:HA	1:C:418:MET:HE1	1.47	0.96
1:A:435:VAL:HG21	1:A:439:PRO:HB3	1.47	0.95
1:B:314:SER:O	4:B:601:HOH:O	1.90	0.90
1:D:435:VAL:CG2	1:D:439:PRO:HB3	2.02	0.90
1:C:387:LEU:HD23	1:C:441:PHE:HE1	1.34	0.89
1:D:435:VAL:HG21	1:D:439:PRO:HB3	1.54	0.89
1:C:379:LYS:HD3	1:C:382:MET:HE3	1.57	0.87
1:D:380:ARG:HA	1:D:444:MET:SD	2.17	0.83
1:C:435:VAL:CG1	1:C:439:PRO:HB3	2.08	0.83
1:D:251:LEU:O	1:D:255:VAL:HG13	1.77	0.82
1:C:387:LEU:HD23	1:C:441:PHE:CE1	2.16	0.81
1:C:305:CYS:HB3	1:C:418:MET:HE2	1.63	0.80
1:D:368:LEU:HD22	1:D:370:VAL:HG23	1.66	0.78
1:D:411:LEU:HD11	1:D:470:LEU:HB3	1.64	0.77
1:A:379:LYS:HG3	4:A:610:HOH:O	1.83	0.77
1:C:259:HIS:HB3	1:C:262:ILE:HD13	1.66	0.77
1:B:216:LYS:HG3	1:B:221:GLU:HG2	1.67	0.76
1:C:379:LYS:HD3	1:C:382:MET:CE	2.16	0.75
1:B:203:THR:HG23	1:B:204:VAL:H	1.52	0.74
1:D:442:GLU:OE2	4:D:601:HOH:O	2.05	0.74
1:C:435:VAL:HG13	1:C:439:PRO:HB3	1.69	0.73
1:A:204:VAL:HA	4:A:632:HOH:O	1.89	0.73
1:B:357:LEU:HB2	1:B:375:ARG:NH1	2.04	0.73
1:D:357:LEU:HD13	1:D:375:ARG:NH1	2.05	0.72
1:C:251:LEU:HD11	1:C:358:ALA:HB3	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:VAL:HG23	1:A:436:PRO:HD2	1.73	0.71
1:D:288:MET:SD	4:D:602:HOH:O	2.48	0.71
1:C:305:CYS:CB	1:C:418:MET:HE2	2.20	0.70
1:D:431:PHE:O	1:D:435:VAL:HG12	1.92	0.70
1:D:234:VAL:HG22	1:D:282:ILE:HD12	1.73	0.70
1:C:492:LYS:O	1:C:496:THR:HG23	1.92	0.70
1:A:431:PHE:O	1:A:435:VAL:HG12	1.91	0.69
1:B:431:PHE:O	1:B:435:VAL:HG22	1.91	0.69
1:D:357:LEU:HD11	1:D:375:ARG:HG2	1.75	0.69
1:C:247:ARG:HD2	1:C:250:GLU:OE2	1.93	0.69
1:C:259:HIS:HB3	1:C:262:ILE:CD1	2.23	0.68
1:D:314:SER:HA	1:D:492:LYS:HD3	1.74	0.68
1:C:392:GLN:HB2	1:C:398:SER:OG	1.92	0.68
1:C:490:ARG:HH21	1:C:490:ARG:HG2	1.58	0.68
1:B:357:LEU:HB2	1:B:375:ARG:HH12	1.57	0.68
1:D:288:MET:HB2	1:D:344:VAL:O	1.92	0.67
1:B:490:ARG:HG3	1:B:490:ARG:HH11	1.60	0.66
1:A:293:ASP:OD2	4:A:602:HOH:O	2.13	0.66
1:D:435:VAL:HG23	1:D:436:PRO:HD2	1.78	0.65
1:B:419:VAL:HA	1:B:424:VAL:HG13	1.78	0.65
1:D:302:THR:HA	1:D:418:MET:HE1	1.77	0.65
1:A:246:PHE:O	1:A:250:GLU:HG2	1.97	0.65
1:B:261:ASN:ND2	1:B:311:SER:HB2	2.12	0.65
1:A:320:HIS:CG	1:A:400:LYS:HD2	2.32	0.65
1:C:462:PHE:HA	1:C:468:THR:OG1	1.96	0.65
1:D:310:LEU:O	1:D:310:LEU:HD23	1.96	0.64
1:A:435:VAL:HG21	1:A:439:PRO:CB	2.22	0.64
1:D:300:LEU:CD2	1:D:308:ILE:HD12	2.27	0.64
1:B:443:ASP:HA	1:B:446:LYS:HE3	1.81	0.63
1:C:302:THR:HA	1:C:418:MET:CE	2.26	0.63
1:C:305:CYS:HB3	1:C:418:MET:CE	2.27	0.62
1:D:251:LEU:HD11	1:D:358:ALA:HB3	1.81	0.62
1:D:305:CYS:CB	1:D:418:MET:HE2	2.30	0.62
1:B:492:LYS:O	1:B:496:THR:HG23	2.01	0.61
1:A:335:ARG:CD	1:A:359:VAL:HG13	2.30	0.61
1:D:300:LEU:HD21	1:D:308:ILE:HD12	1.81	0.61
1:C:458:PRO:HG2	1:C:461:TRP:CD1	2.35	0.61
1:B:388:ASP:OD1	1:B:390:THR:HB	2.00	0.60
1:C:412:TRP:CZ2	1:C:416:ARG:HD2	2.36	0.60
1:C:247:ARG:HA	1:C:250:GLU:HG2	1.84	0.60
1:B:442:GLU:O	1:B:446:LYS:HG2	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:314:SER:OG	1:D:492:LYS:HE3	2.02	0.60
1:B:462:PHE:HA	1:B:468:THR:CG2	2.32	0.60
1:A:255:VAL:O	1:A:257:LEU:N	2.35	0.59
1:C:385:GLU:HG3	1:C:391:ILE:HD12	1.84	0.59
1:D:251:LEU:HA	1:D:255:VAL:CG1	2.33	0.59
1:D:330:PRO:HG3	1:D:360:MET:CE	2.32	0.59
1:D:453:GLN:O	1:D:454:ARG:HD2	2.02	0.59
1:B:255:VAL:O	1:B:257:LEU:N	2.35	0.58
1:C:368:LEU:CD1	1:C:396:PHE:HB2	2.34	0.57
1:C:292:TYR:O	1:C:296:GLN:HG2	2.03	0.57
1:A:332:ILE:HG12	1:A:360:MET:HG2	1.86	0.57
1:C:368:LEU:HD12	1:C:396:PHE:HB2	1.87	0.57
1:C:458:PRO:HG2	1:C:461:TRP:NE1	2.20	0.57
1:C:379:LYS:HA	1:C:382:MET:HE3	1.87	0.56
1:C:334:HIS:O	1:C:335:ARG:HB2	2.05	0.56
1:C:420:SER:O	1:C:423:ILE:HG22	2.06	0.56
1:B:251:LEU:O	1:B:255:VAL:O	2.24	0.56
1:B:452:GLN:NE2	1:B:480:GLN:OE1	2.37	0.56
1:D:457:ILE:HD11	1:D:478:TRP:HZ3	1.70	0.55
1:D:330:PRO:HG3	1:D:360:MET:HE2	1.88	0.55
1:B:489:LEU:HD11	1:B:493:LYS:HE2	1.88	0.55
1:D:368:LEU:HD23	1:D:369:ASP:N	2.21	0.55
1:A:417:ARG:NH1	1:A:426:ASP:HA	2.22	0.55
1:B:453:GLN:O	1:B:454:ARG:HD2	2.07	0.55
1:C:288:MET:HE1	1:C:346:LYS:HA	1.89	0.54
1:A:271:THR:HG22	1:A:278:GLN:HB2	1.88	0.54
1:A:273:ARG:HG2	1:A:274:HIS:N	2.23	0.54
1:A:270:MET:HG2	1:A:271:THR:N	2.20	0.54
1:B:334:HIS:O	1:B:335:ARG:HB2	2.08	0.53
1:A:298:THR:HG22	1:A:299:THR:O	2.07	0.53
1:A:426:ASP:OD1	1:A:426:ASP:N	2.41	0.53
1:C:435:VAL:HG11	1:C:439:PRO:HB3	1.90	0.53
1:C:490:ARG:NH2	1:C:490:ARG:HG2	2.21	0.53
1:D:334:HIS:O	1:D:335:ARG:HB2	2.09	0.53
1:A:496:THR:O	1:B:307:ARG:NH1	2.42	0.53
1:A:251:LEU:HD11	1:A:358:ALA:HB3	1.91	0.52
1:C:360:MET:O	1:C:368:LEU:HA	2.09	0.52
1:C:401:ARG:NH1	4:C:601:HOH:O	2.42	0.52
1:D:412:TRP:CZ2	1:D:416:ARG:HD2	2.44	0.52
1:B:419:VAL:HA	1:B:424:VAL:CG1	2.40	0.52
1:D:435:VAL:HG21	1:D:439:PRO:CB	2.32	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:MET:HB2	4:A:612:HOH:O	2.09	0.52
1:C:245:TRP:NE1	1:C:268:SER:OG	2.42	0.52
1:A:208:ILE:CD1	1:A:282:ILE:HD13	2.40	0.52
1:A:402:VAL:HG13	4:A:614:HOH:O	2.09	0.52
1:B:420:SER:O	1:B:423:ILE:HG22	2.09	0.52
1:D:476:GLU:HB2	1:D:486:LEU:HD11	1.91	0.52
1:B:448:VAL:O	1:B:452:GLN:HA	2.10	0.51
1:B:477:CYS:O	1:B:485:ARG:HD3	2.11	0.51
1:C:288:MET:CE	1:C:346:LYS:HA	2.40	0.51
1:C:318:HIS:O	1:C:329:LYS:HE2	2.10	0.51
1:D:455:PRO:O	1:D:457:ILE:HD12	2.10	0.51
1:D:357:LEU:HD13	1:D:375:ARG:CZ	2.40	0.51
1:C:489:LEU:HD23	1:C:489:LEU:O	2.10	0.51
1:A:412:TRP:CZ2	1:A:416:ARG:HD2	2.46	0.51
1:B:416:ARG:NH1	1:B:426:ASP:O	2.44	0.51
1:A:204:VAL:N	4:A:605:HOH:O	2.43	0.51
1:B:379:LYS:HA	1:B:382:MET:CE	2.41	0.50
1:A:235:LYS:HE2	1:A:237:PHE:CZ	2.47	0.50
1:B:317:ALA:O	1:B:321:ILE:HG12	2.10	0.50
1:B:490:ARG:HG3	1:B:490:ARG:NH1	2.23	0.50
1:C:435:VAL:HG13	1:C:436:PRO:HD2	1.92	0.50
1:D:300:LEU:HD21	1:D:308:ILE:CD1	2.41	0.50
1:D:301:ASP:OD2	1:D:303:VAL:HG12	2.12	0.50
1:A:292:TYR:O	1:A:296:GLN:HG2	2.12	0.50
1:B:443:ASP:HA	1:B:446:LYS:CE	2.41	0.50
1:B:458:PRO:HG2	1:B:461:TRP:NE1	2.27	0.50
1:D:249:THR:HG22	1:D:253:ASN:ND2	2.27	0.50
1:C:435:VAL:HG23	1:C:447:VAL:HG21	1.94	0.49
1:D:357:LEU:HD11	1:D:375:ARG:CG	2.42	0.49
1:D:305:CYS:HB3	1:D:418:MET:CE	2.42	0.49
1:A:251:LEU:O	1:A:255:VAL:O	2.31	0.49
1:A:214:VAL:HG12	2:A:501:B4E:N17	2.28	0.49
1:C:246:PHE:O	1:C:250:GLU:HG2	2.12	0.49
1:C:237:PHE:HB2	1:C:279:LEU:HB2	1.94	0.49
1:A:235:LYS:HE2	1:A:237:PHE:CE2	2.47	0.49
1:D:462:PHE:HA	1:D:468:THR:OG1	2.13	0.49
1:C:332:ILE:HG12	1:C:360:MET:HG2	1.95	0.49
1:D:255:VAL:O	1:D:257:LEU:N	2.44	0.48
1:D:332:ILE:HG12	1:D:360:MET:HG2	1.94	0.48
1:C:411:LEU:HD23	1:C:474:MET:HE2	1.94	0.48
1:D:385:GLU:HG3	1:D:391:ILE:HD12	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:458:PRO:HG2	1:D:461:TRP:CD1	2.48	0.48
1:A:248:GLU:HG3	1:A:355:LEU:HB2	1.95	0.48
1:C:411:LEU:HD22	1:C:474:MET:HG3	1.95	0.48
1:D:251:LEU:HD23	1:D:255:VAL:HG11	1.96	0.48
1:D:472:LYS:O	1:D:476:GLU:HG3	2.13	0.48
1:D:384:PRO:HB2	1:D:480:GLN:OE1	2.14	0.48
1:A:379:LYS:HA	1:A:382:MET:CE	2.44	0.48
1:B:462:PHE:HA	1:B:468:THR:HG22	1.96	0.48
1:A:334:HIS:O	1:A:335:ARG:HB2	2.14	0.47
1:C:286:HIS:CE1	1:C:345:LYS:HG2	2.49	0.47
1:C:299:THR:HG23	1:C:419:VAL:CG2	2.45	0.47
1:B:286:HIS:CE1	1:B:345:LYS:HG2	2.50	0.47
1:C:300:LEU:O	1:C:419:VAL:HG23	2.15	0.47
1:D:343:LEU:O	1:D:350:CYS:HA	2.15	0.47
1:A:357:LEU:HB2	4:A:609:HOH:O	2.16	0.46
1:B:270:MET:HG2	1:B:271:THR:N	2.22	0.46
1:B:299:THR:HG22	1:B:417:ARG:NH2	2.30	0.46
1:C:379:LYS:HA	1:C:382:MET:CE	2.45	0.46
1:C:299:THR:CG2	1:C:419:VAL:CG2	2.93	0.46
1:C:363:GLN:HA	4:C:604:HOH:O	2.15	0.46
1:D:296:GLN:NE2	1:D:427:TYR:CD2	2.83	0.46
1:D:457:ILE:N	1:D:457:ILE:HD12	2.31	0.46
1:B:261:ASN:O	1:B:351:CYS:HA	2.16	0.46
1:C:411:LEU:HD23	1:C:474:MET:CE	2.45	0.46
1:D:411:LEU:HD12	1:D:474:MET:HG3	1.97	0.45
1:D:458:PRO:HG2	1:D:461:TRP:NE1	2.32	0.45
1:B:458:PRO:HG2	1:B:461:TRP:CE2	2.52	0.45
1:C:384:PRO:HD3	1:C:405:TRP:CZ2	2.51	0.45
1:C:377:GLY:O	1:C:382:MET:HE2	2.15	0.45
1:B:335:ARG:HD3	1:B:357:LEU:O	2.17	0.45
1:A:435:VAL:HG22	1:A:436:PRO:O	2.17	0.45
1:C:251:LEU:O	1:C:255:VAL:O	2.34	0.45
1:C:411:LEU:CD2	1:C:474:MET:HG3	2.47	0.45
1:A:214:VAL:O	1:A:214:VAL:CG1	2.64	0.45
1:A:333:ALA:HB3	1:A:359:VAL:HG22	1.98	0.45
1:D:305:CYS:CB	1:D:418:MET:CE	2.95	0.45
1:B:247:ARG:NH1	1:B:360:MET:SD	2.90	0.45
1:C:288:MET:HE1	1:C:345:LYS:O	2.17	0.45
1:C:476:GLU:HB3	1:C:486:LEU:HG	1.99	0.45
1:C:251:LEU:O	1:C:255:VAL:HG13	2.17	0.44
1:C:285:TYR:CE2	1:C:287:GLU:HA	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:288:MET:HA	4:D:602:HOH:O	2.17	0.44
1:D:305:CYS:HB3	1:D:418:MET:HE2	1.98	0.44
1:D:435:VAL:CG2	1:D:436:PRO:HD2	2.47	0.44
1:D:411:LEU:HD12	1:D:474:MET:HE3	2.00	0.44
1:A:208:ILE:HD12	1:A:282:ILE:HD13	1.98	0.44
1:C:264:GLY:O	1:C:283:THR:HB	2.18	0.44
1:D:298:THR:HG22	1:D:299:THR:O	2.17	0.44
1:C:270:MET:HE2	1:C:279:LEU:HD21	1.99	0.43
1:B:207:ASP:HB2	4:B:602:HOH:O	2.18	0.43
1:C:299:THR:HG23	1:C:419:VAL:HG23	1.99	0.43
1:D:222:VAL:HA	1:D:234:VAL:O	2.19	0.43
1:A:335:ARG:HD3	1:A:359:VAL:HG13	2.00	0.43
1:A:379:LYS:HA	1:A:382:MET:HG3	2.00	0.43
1:D:334:HIS:N	1:D:403:ASP:OD2	2.45	0.43
1:D:435:VAL:HG22	1:D:439:PRO:HB3	1.93	0.43
1:D:318:HIS:O	1:D:329:LYS:HE2	2.18	0.43
1:B:379:LYS:HA	1:B:382:MET:HE2	2.01	0.43
1:C:379:LYS:CD	1:C:382:MET:HE3	2.39	0.43
1:A:333:ALA:O	1:A:358:ALA:HA	2.19	0.42
1:B:211:LEU:HD11	1:B:226:SER:HB2	2.01	0.42
1:C:333:ALA:O	1:C:358:ALA:HA	2.19	0.42
1:D:251:LEU:HD11	1:D:358:ALA:CB	2.49	0.42
1:C:385:GLU:HG3	1:C:391:ILE:CD1	2.48	0.42
1:C:428:LYS:HD3	1:C:428:LYS:HA	1.83	0.42
1:C:251:LEU:HD21	1:C:360:MET:SD	2.60	0.42
1:D:286:HIS:CE1	1:D:345:LYS:HG2	2.55	0.42
2:D:501:B4E:N15	2:D:501:B4E:C19	2.83	0.42
1:A:271:THR:CG2	1:A:278:GLN:HB2	2.49	0.42
1:C:461:TRP:HB3	1:C:467:LEU:HB3	2.00	0.42
1:C:305:CYS:CB	1:C:418:MET:CE	2.93	0.42
1:C:387:LEU:CD1	1:C:448:VAL:HB	2.49	0.42
1:B:443:ASP:HA	1:B:446:LYS:NZ	2.35	0.42
1:C:204:VAL:HG23	1:C:205:ALA:N	2.34	0.42
1:C:461:TRP:HZ3	1:C:474:MET:HE3	1.85	0.42
1:A:251:LEU:HD21	1:A:360:MET:SD	2.60	0.41
1:A:306:LEU:HD23	1:A:306:LEU:HA	1.88	0.41
1:B:427:TYR:C	1:B:427:TYR:CD1	2.93	0.41
1:B:412:TRP:CZ2	1:B:416:ARG:HD3	2.55	0.41
1:C:428:LYS:HB3	1:C:432:TYR:CG	2.54	0.41
1:D:400:LYS:O	1:D:404:ILE:HG12	2.21	0.41
1:D:331:ALA:O	1:D:360:MET:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:411:LEU:CD2	1:C:474:MET:HE2	2.49	0.41
1:D:330:PRO:CG	1:D:360:MET:HE3	2.51	0.41
1:A:273:ARG:HG2	1:A:274:HIS:H	1.85	0.41
1:A:214:VAL:CG1	2:A:501:B4E:N17	2.84	0.41
1:B:214:VAL:HG13	2:B:501:B4E:N17	2.36	0.41
1:B:376:VAL:HG22	1:B:377:GLY:N	2.36	0.41
1:A:286:HIS:HD2	4:A:626:HOH:O	2.03	0.41
1:C:261:ASN:ND2	1:C:311:SER:HB2	2.36	0.41
1:C:301:ASP:OD2	1:C:303:VAL:HG12	2.21	0.41
1:B:397:ASP:OD1	1:B:400:LYS:NZ	2.33	0.41
1:C:204:VAL:HG23	1:C:205:ALA:H	1.86	0.41
1:C:255:VAL:HG13	1:C:255:VAL:O	2.21	0.41
1:C:233:ALA:O	1:C:282:ILE:HA	2.21	0.41
1:A:376:VAL:HG22	1:A:377:GLY:N	2.36	0.41
1:A:457:ILE:HG21	1:A:462:PHE:CZ	2.57	0.40
1:D:448:VAL:O	1:D:452:GLN:HA	2.21	0.40
1:A:306:LEU:HA	1:A:309:VAL:HG22	2.03	0.40
1:C:379:LYS:HA	1:C:382:MET:HG3	2.03	0.40
1:C:435:VAL:HG12	1:C:436:PRO:O	2.21	0.40
1:C:418:MET:HG2	1:C:419:VAL:N	2.36	0.40
1:D:251:LEU:O	1:D:255:VAL:O	2.39	0.40
1:D:309:VAL:HG11	1:D:411:LEU:HD22	2.03	0.40
1:D:342:ILE:CG2	1:D:350:CYS:HB3	2.52	0.40

There are no symmetry-related clashes.

### 5.3 Torsion angles [\(i\)](#)

#### 5.3.1 Protein backbone [\(i\)](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	287/301 (95%)	279 (97%)	8 (3%)	0	100 100
1	B	288/301 (96%)	276 (96%)	12 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	287/301 (95%)	278 (97%)	9 (3%)	0	100	100
1	D	288/301 (96%)	281 (98%)	7 (2%)	0	100	100
All	All	1150/1204 (96%)	1114 (97%)	36 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	241/270 (89%)	235 (98%)	6 (2%)	47	61
1	B	239/270 (88%)	232 (97%)	7 (3%)	42	55
1	C	219/270 (81%)	213 (97%)	6 (3%)	44	58
1	D	204/270 (76%)	194 (95%)	10 (5%)	25	34
All	All	903/1080 (84%)	874 (97%)	29 (3%)	39	51

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	204	VAL
1	A	209	THR
1	A	326	THR
1	A	379	LYS
1	A	419	VAL
1	A	426	ASP
1	B	203	THR
1	B	341	ASN
1	B	390	THR
1	B	416	ARG
1	B	424	VAL
1	B	446	LYS
1	B	460	ARG
1	C	299	THR

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Mol	Chain	Res	Type
1	C	368	LEU
1	C	411	LEU
1	C	426	ASP
1	C	441	PHE
1	C	476	GLU
1	D	204	VAL
1	D	222	VAL
1	D	255	VAL
1	D	277	THR
1	D	310	LEU
1	D	357	LEU
1	D	368	LEU
1	D	426	ASP
1	D	444	MET
1	D	460	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [\(i\)](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [\(i\)](#)

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

### 5.5 Carbohydrates [\(i\)](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [\(i\)](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	B4E	D	501	-	20,26,26	1.76	2 (10%)	23,36,36	1.82	3 (13%)
2	B4E	C	501	-	20,26,26	1.86	2 (10%)	23,36,36	1.93	2 (8%)
2	B4E	A	501	-	20,26,26	1.78	3 (15%)	23,36,36	1.65	3 (13%)
3	EDO	A	502	-	3,3,3	0.46	0	2,2,2	0.32	0
2	B4E	B	501	-	20,26,26	1.88	2 (10%)	23,36,36	1.91	4 (17%)
3	EDO	A	503	-	3,3,3	0.48	0	2,2,2	0.29	0

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	B4E	D	501	-	-	9/14/14/14	0/3/4/4
2	B4E	C	501	-	-	4/14/14/14	0/3/4/4
2	B4E	A	501	-	-	6/14/14/14	0/3/4/4
3	EDO	A	502	-	-	0/1/1/1	-
2	B4E	B	501	-	-	5/14/14/14	0/3/4/4
3	EDO	A	503	-	-	0/1/1/1	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	B4E	C08-C06	-6.27	1.38	1.49
2	D	501	B4E	C08-C06	-5.64	1.39	1.49
2	C	501	B4E	C08-C06	-5.50	1.40	1.49
2	A	501	B4E	C08-C06	-5.40	1.40	1.49
2	C	501	B4E	C14-N17	5.08	1.43	1.36
2	B	501	B4E	C14-N17	4.54	1.43	1.36
2	A	501	B4E	C14-N17	4.54	1.43	1.36
2	D	501	B4E	C14-N17	4.34	1.42	1.36
2	A	501	B4E	O23-C04	2.20	1.42	1.37

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	B4E	C09-N10-N11	7.86	109.42	103.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	B4E	C09-N10-N11	7.19	108.93	103.70
2	D	501	B4E	C09-N10-N11	6.39	108.35	103.70
2	A	501	B4E	C09-N10-N11	5.80	107.92	103.70
2	D	501	B4E	C14-N15-C16	3.42	120.71	117.59
2	C	501	B4E	C14-N15-C16	3.16	120.46	117.59
2	B	501	B4E	C03-C02-CL1	2.64	122.45	119.15
2	D	501	B4E	C18-N17-C14	-2.61	119.25	123.38
2	A	501	B4E	C18-N17-C14	-2.33	119.70	123.38
2	A	501	B4E	C03-C02-CL1	2.27	121.98	119.15
2	B	501	B4E	C18-N17-C14	-2.24	119.83	123.38
2	B	501	B4E	C14-N15-C16	2.13	119.53	117.59

There are no chirality outliers.

All (24) torsion outliers are listed below:

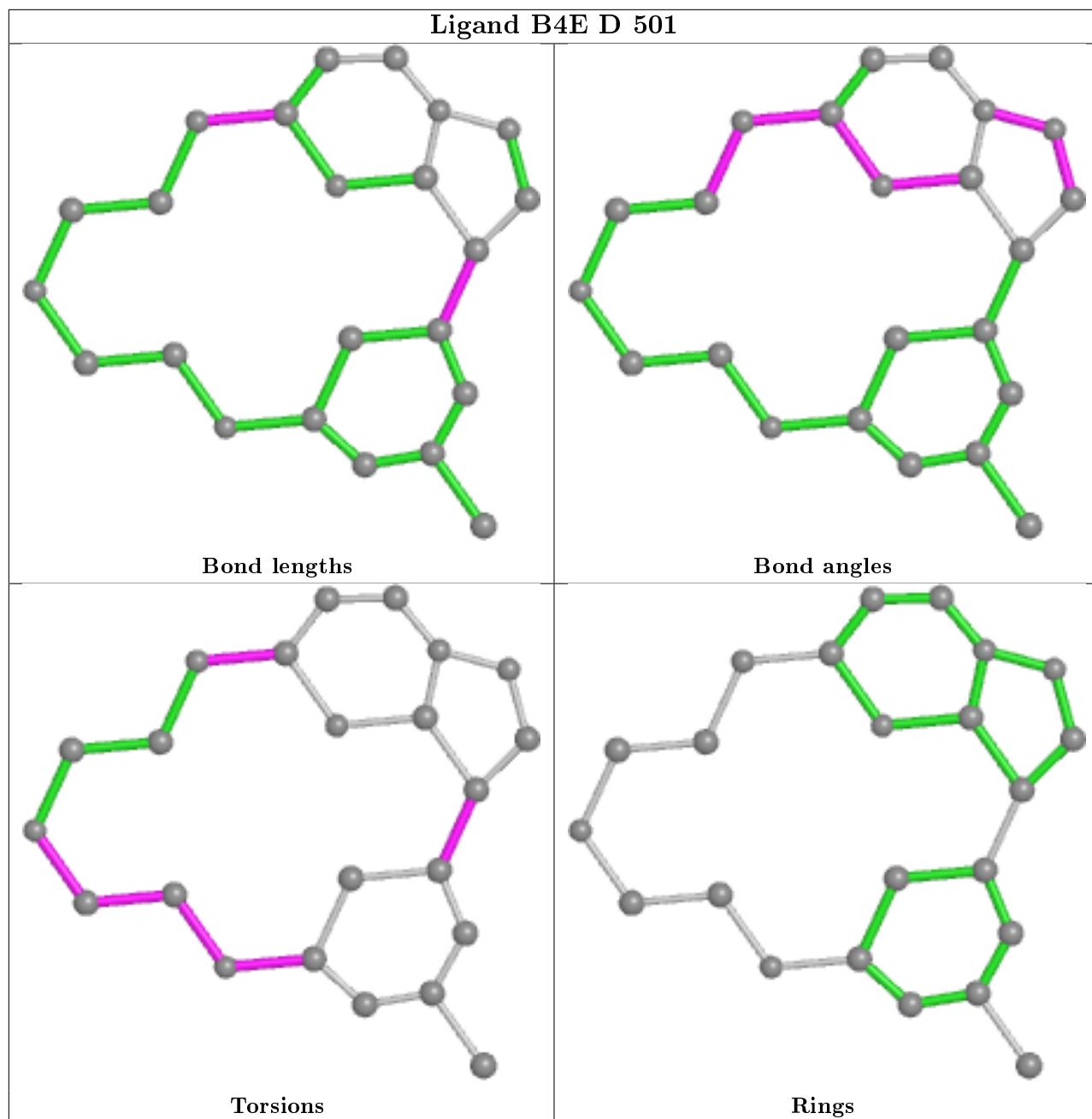
Mol	Chain	Res	Type	Atoms
2	D	501	B4E	C05-C06-C08-C16
2	D	501	B4E	C07-C06-C08-C16
2	D	501	B4E	C13-C14-N17-C18
2	D	501	B4E	N15-C14-N17-C18
2	D	501	B4E	C22-C21-O20-C19
2	A	501	B4E	O20-C21-C22-O23
2	B	501	B4E	O20-C21-C22-O23
2	A	501	B4E	C03-C04-O23-C22
2	B	501	B4E	C03-C04-O23-C22
2	A	501	B4E	C05-C04-O23-C22
2	B	501	B4E	C05-C04-O23-C22
2	D	501	B4E	O20-C21-C22-O23
2	A	501	B4E	N17-C18-C19-O20
2	D	501	B4E	C03-C04-O23-C22
2	D	501	B4E	C05-C04-O23-C22
2	D	501	B4E	C21-C22-O23-C04
2	B	501	B4E	C21-C22-O23-C04
2	B	501	B4E	C22-C21-O20-C19
2	C	501	B4E	C05-C06-C08-C16
2	A	501	B4E	C21-C22-O23-C04
2	C	501	B4E	C21-C22-O23-C04
2	C	501	B4E	C07-C06-C08-C16
2	A	501	B4E	C07-C06-C08-C16
2	C	501	B4E	O20-C21-C22-O23

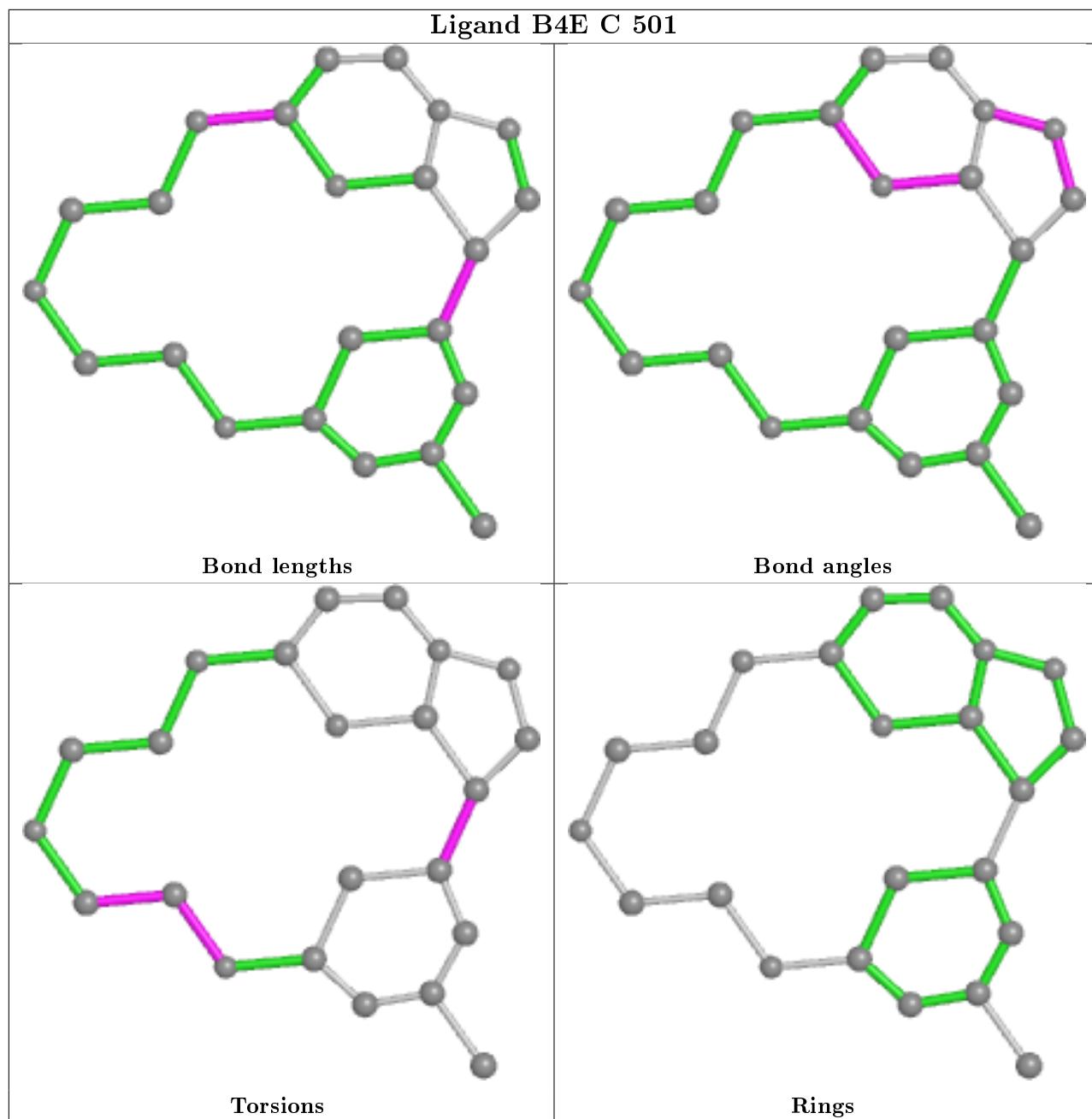
There are no ring outliers.

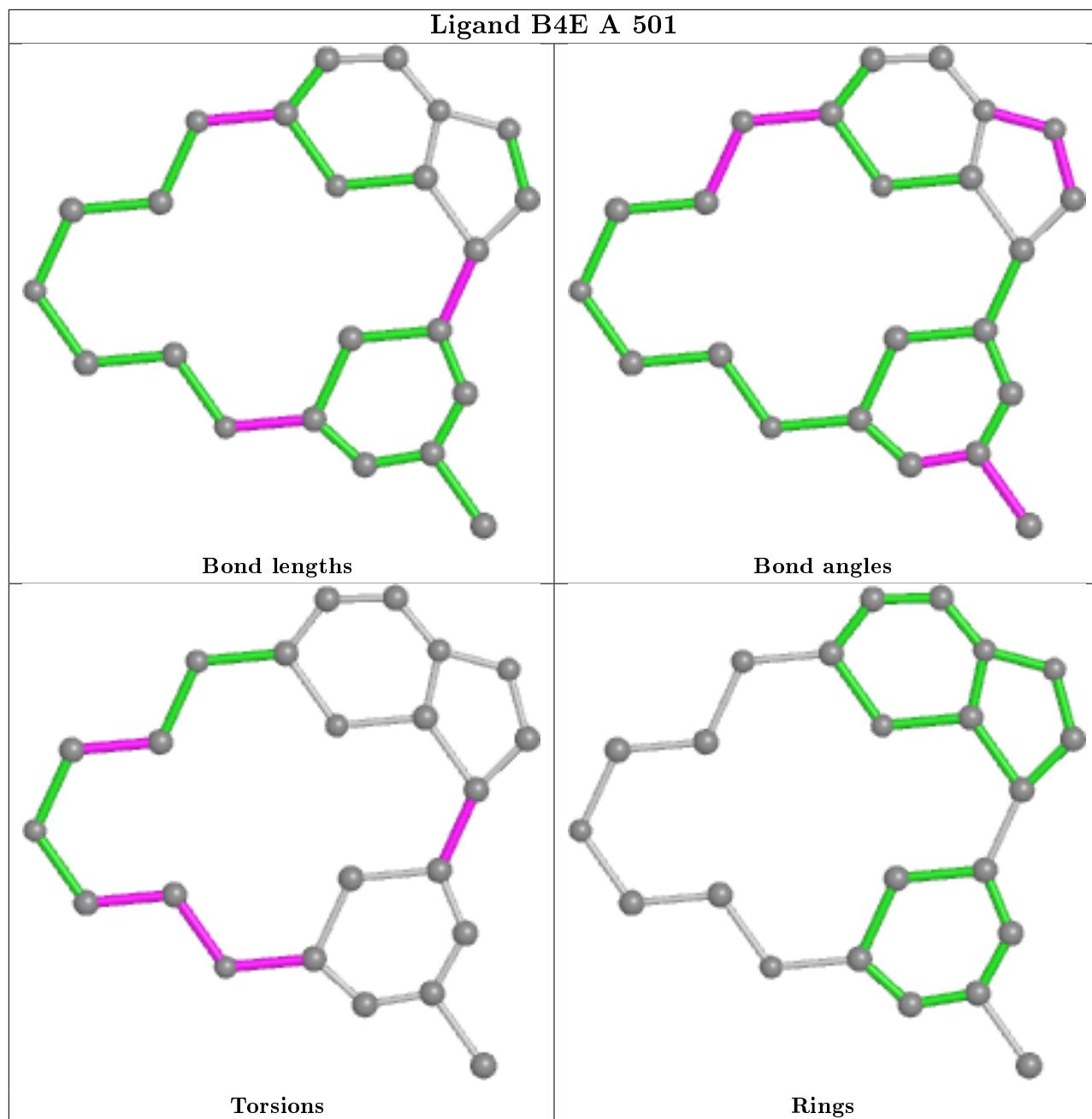
3 monomers are involved in 4 short contacts:

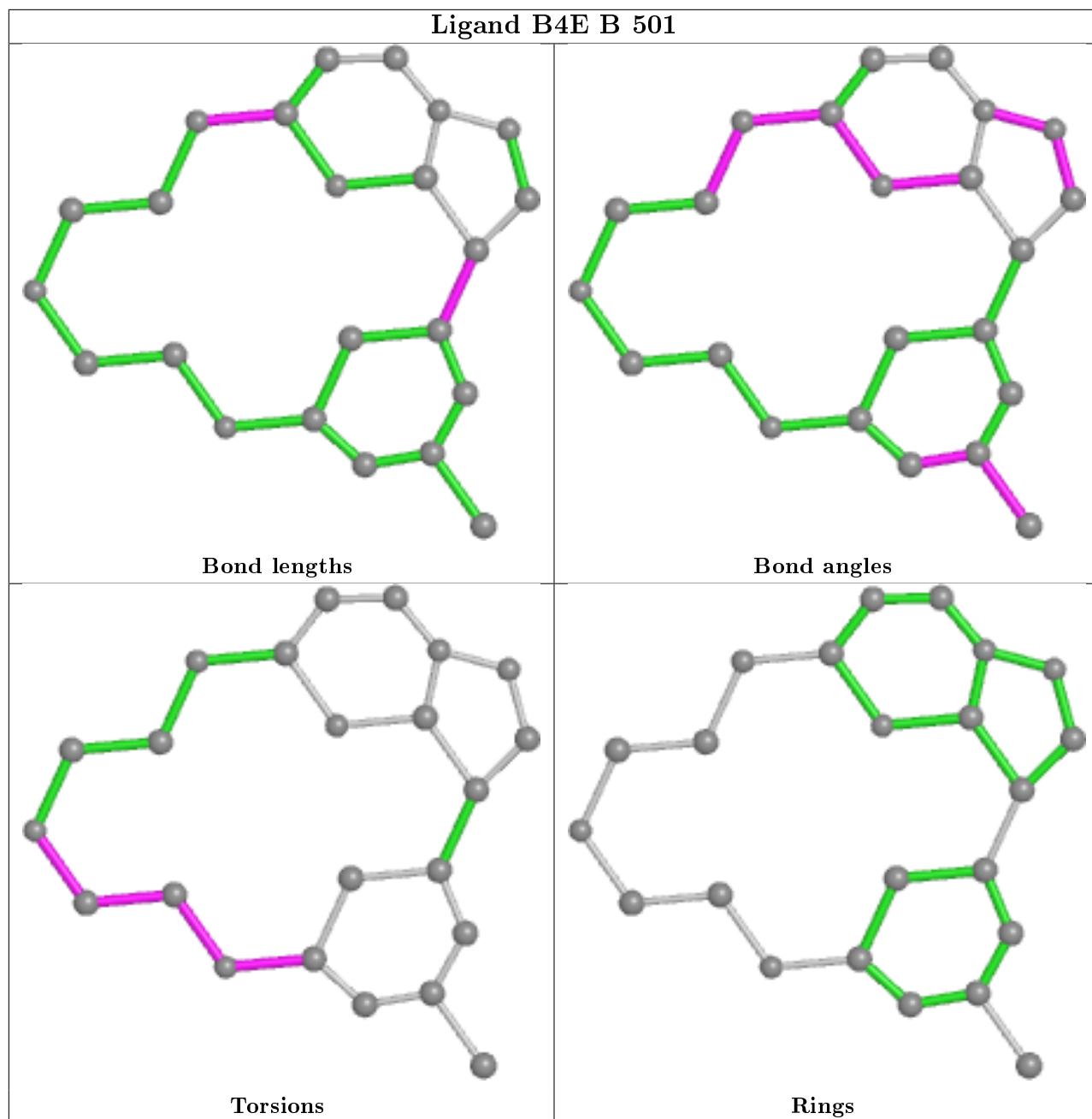
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	B4E	1	0
2	A	501	B4E	2	0
2	B	501	B4E	1	0

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 than 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.









## 5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

## 6 Fit of model and data i

### 6.1 Protein, DNA and RNA chains i

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	A	291/301 (96%)	0.07	2 (0%) 87 91	48, 74, 106, 130	0
1	B	292/301 (97%)	0.18	7 (2%) 59 67	53, 85, 118, 141	0
1	C	291/301 (96%)	0.24	9 (3%) 49 58	58, 90, 127, 144	0
1	D	292/301 (97%)	0.42	24 (8%) 11 16	30, 102, 134, 166	0
All	All	1166/1204 (96%)	0.23	42 (3%) 42 51	30, 88, 126, 166	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	324	PHE	5.0
1	D	355	LEU	4.2
1	D	326	THR	4.2
1	D	431	PHE	4.1
1	D	324	PHE	4.0
1	D	325	GLY	4.0
1	D	219	TYR	3.4
1	B	473	LEU	3.3
1	D	381	TYR	3.1
1	D	488	ALA	3.1
1	D	337	LEU	3.1
1	B	461	TRP	2.9
1	C	265	PHE	2.9
1	B	483	SER	2.9
1	D	214	VAL	2.8
1	D	427	TYR	2.7
1	D	295	LEU	2.7
1	B	498	ILE	2.6
1	C	363	GLN	2.5
1	C	270	MET	2.5
1	B	297	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	291	LEU	2.4
1	D	251	LEU	2.4
1	C	204	VAL	2.4
1	D	353	ALA	2.4
1	D	368	LEU	2.4
1	D	342	ILE	2.3
1	D	396	PHE	2.3
1	D	407	PHE	2.3
1	D	478	TRP	2.3
1	B	470	LEU	2.2
1	C	263	LEU	2.2
1	C	370	VAL	2.1
1	B	465	PRO	2.1
1	D	289	GLY	2.1
1	D	430	PRO	2.1
1	C	210	LEU	2.1
1	D	470	LEU	2.1
1	C	205	ALA	2.1
1	A	411	LEU	2.1
1	D	270	MET	2.0
1	A	275	SER	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [\(i\)](#)

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

## 6.3 Carbohydrates [\(i\)](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [\(i\)](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

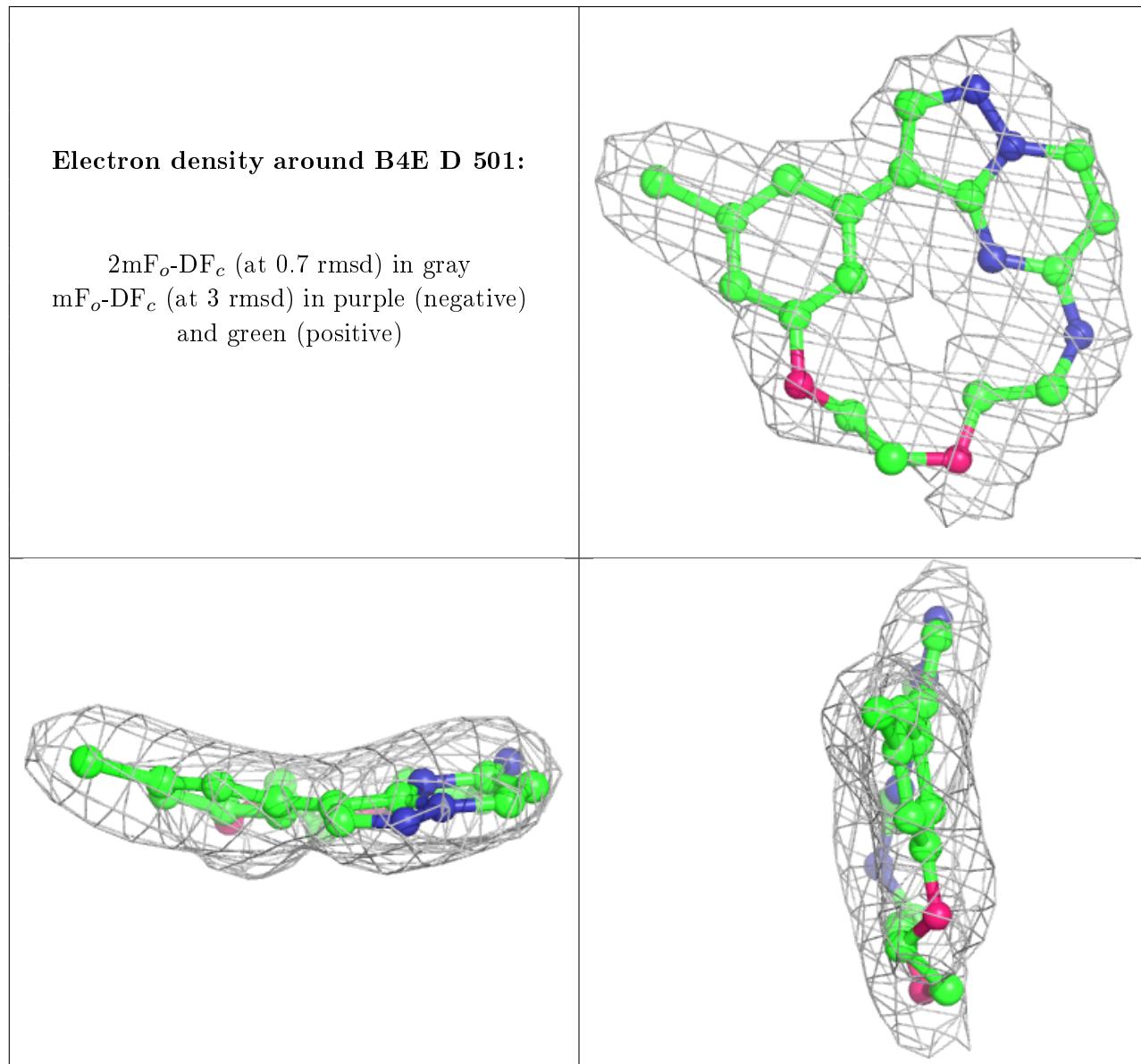
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	EDO	A	503	4/4	0.66	0.18	101,121,121,121	0
3	EDO	A	502	4/4	0.81	0.12	108,130,131,131	0

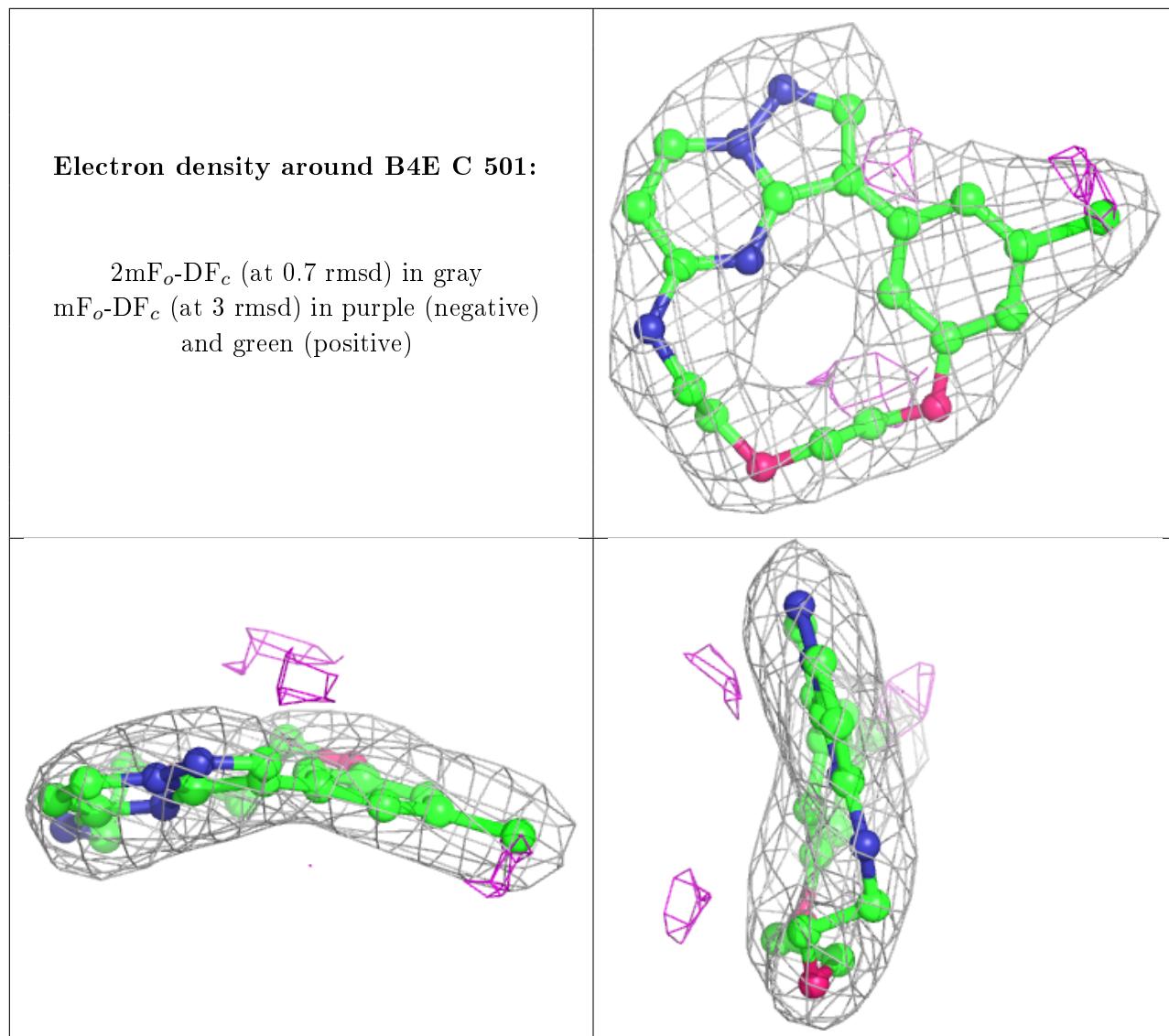
*Continued on next page...*

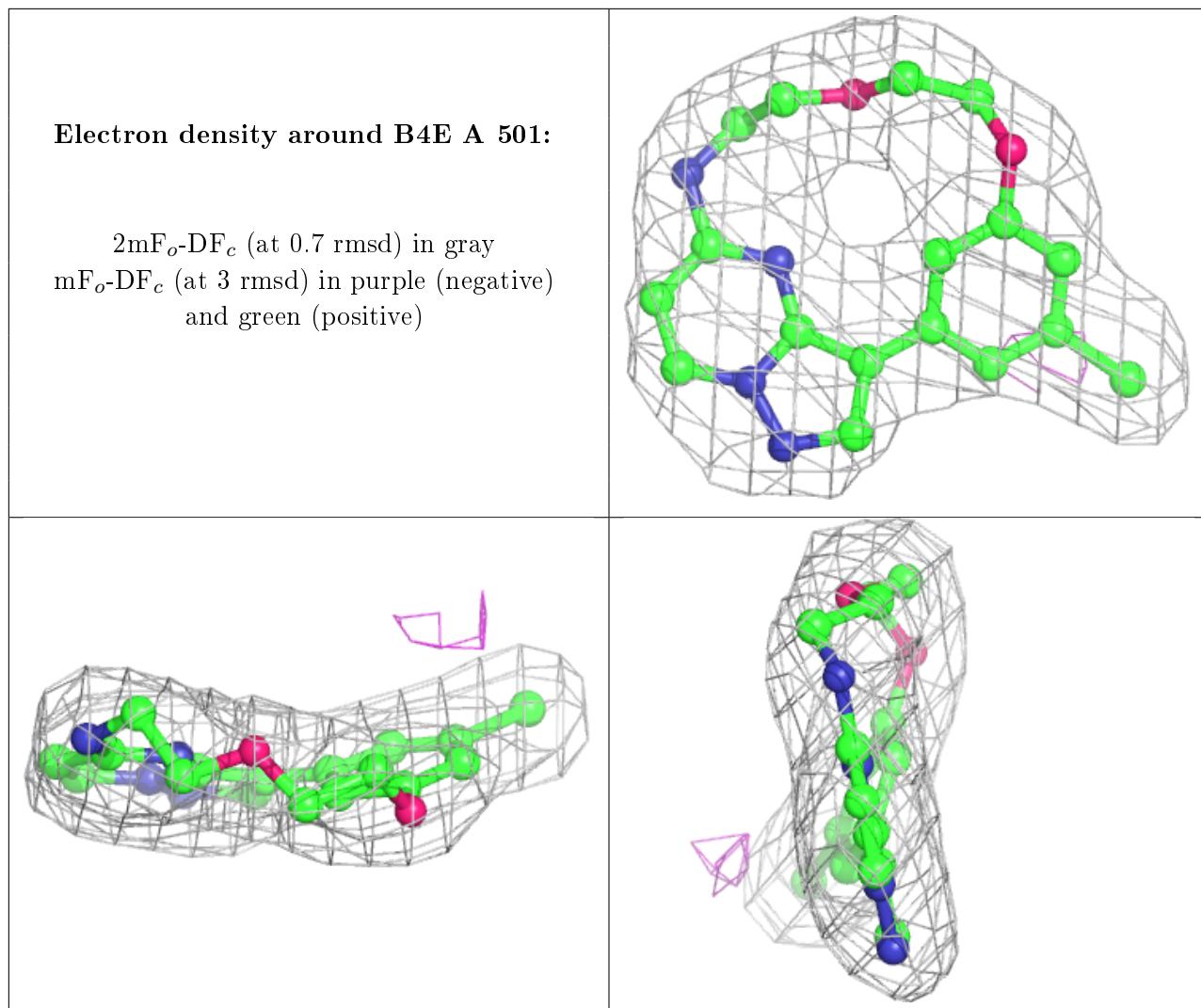
*Continued from previous page...*

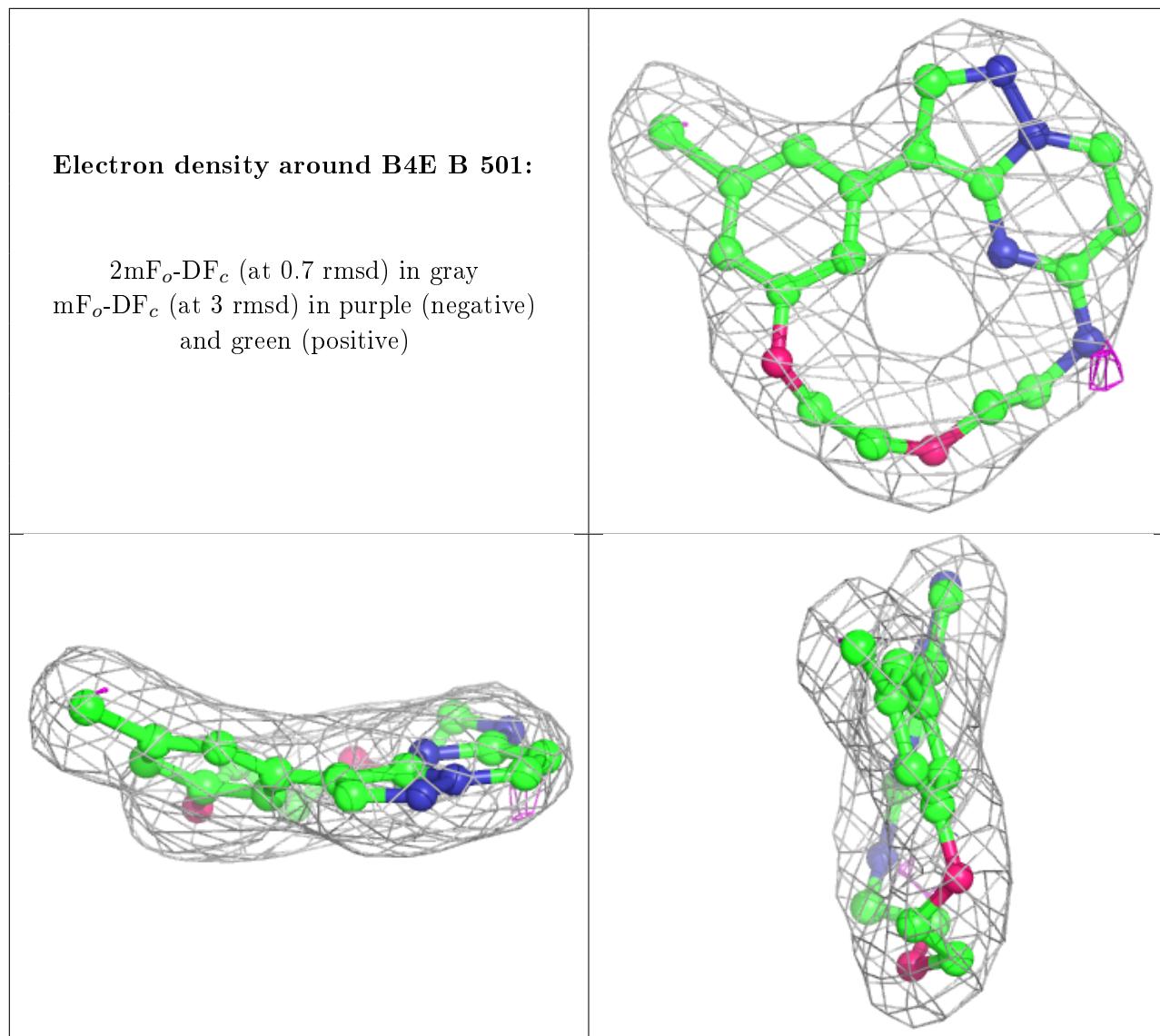
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	B4E	D	501	23/23	0.90	0.18	76,78,81,82	0
2	B4E	C	501	23/23	0.93	0.18	74,75,76,78	0
2	B4E	A	501	23/23	0.96	0.17	57,61,64,65	0
2	B4E	B	501	23/23	0.97	0.20	59,62,67,68	0

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.









## 6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.