



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:22 AM UTC

PDB ID : 2IDK / pdb_00002idk
Title : Crystal Structure of Rat Glycine N-Methyltransferase Complexed With Folate
Authors : Luka, Z.; Pakhomova, S.; Loukachevitch, L.V.; Egli, M.; Newcomer, M.E.; Wagner, C.
Deposited on : 2006-09-15
Resolution : 2.55 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

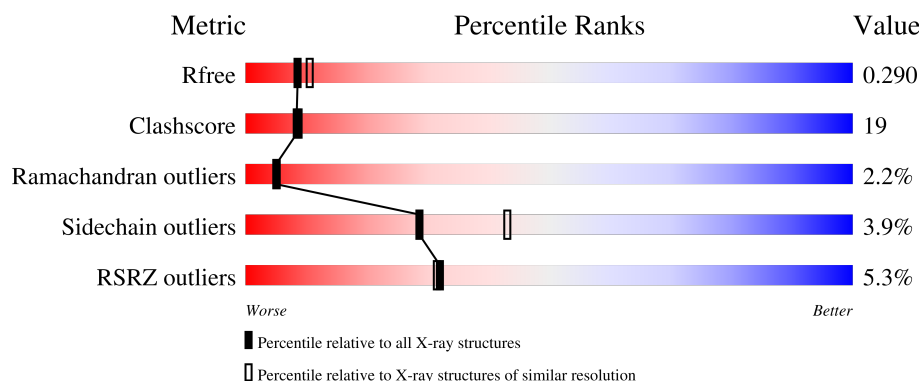
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	292	5% 57% 36% • 5%
1	B	292	5% 60% 34% • •
1	C	292	5% 56% 38% • •
1	D	292	4% 61% 33% • •

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	C2F	A	1410	-	-	X	-
2	C2F	B	1420	-	-	X	-

2 Entry composition [i](#)

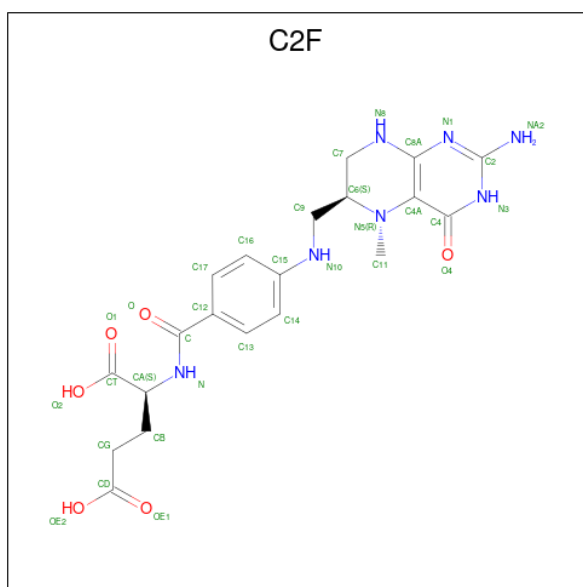
There are 3 unique types of molecules in this entry. The entry contains 8848 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycine N-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	277	Total	C	N	O	S	0	0	0
			2140	1361	372	396	11			
1	B	285	Total	C	N	O	S	0	0	0
			2219	1410	385	413	11			
1	C	281	Total	C	N	O	S	0	0	0
			2176	1385	379	401	11			
1	D	287	Total	C	N	O	S	0	0	0
			2233	1419	386	417	11			

- Molecule 2 is 5-METHYL-5,6,7,8-TETRAHYDROFOLIC ACID (CCD ID: C2F) (formula: $C_{20}H_{25}N_7O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			33	20	7	6		
2	B	1	Total	C	N	O	0	0
			33	20	7	6		

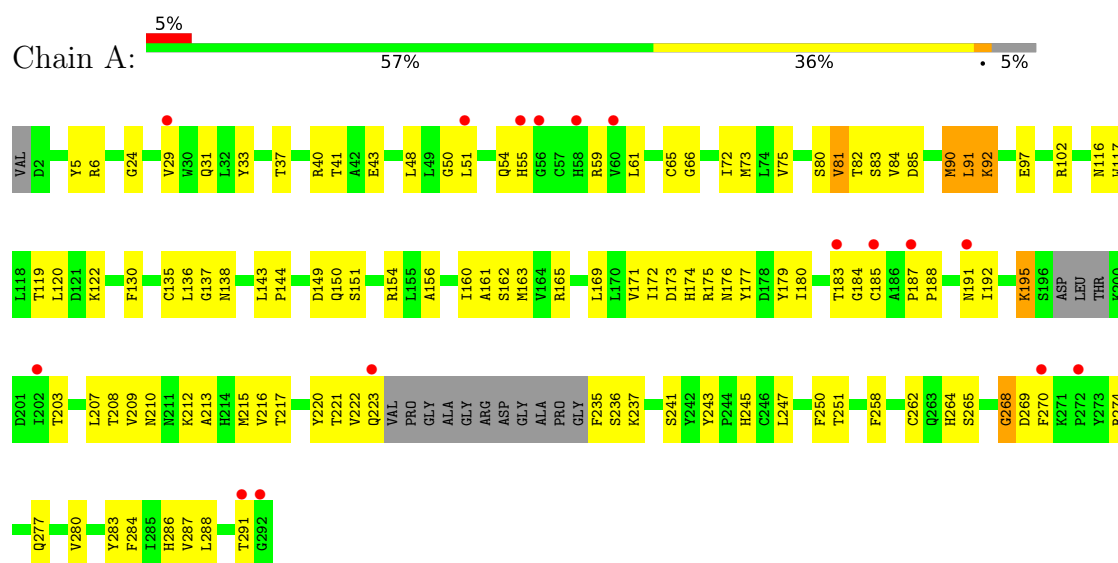
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	O 1	0	0
3	B	5	Total 5	O 5	0	0
3	C	4	Total 4	O 4	0	0
3	D	4	Total 4	O 4	0	0

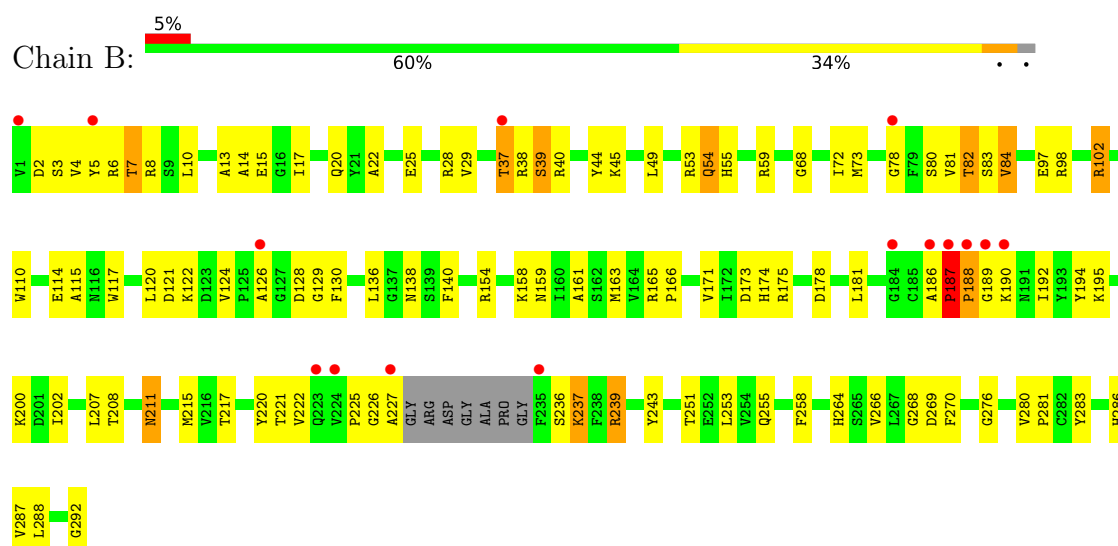
3 Residue-property plots

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: Glycine N-methyltransferase

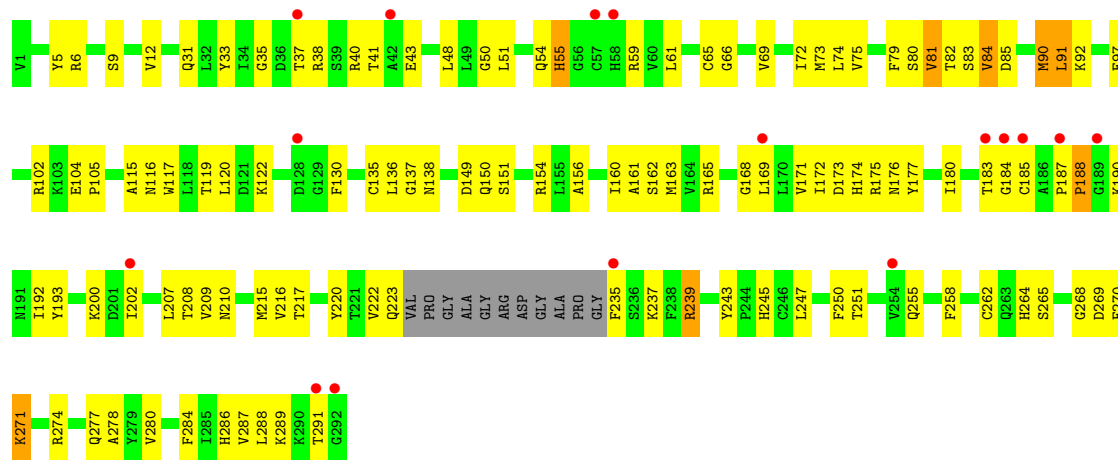


• Molecule 1: Glycine N-methyltransferase

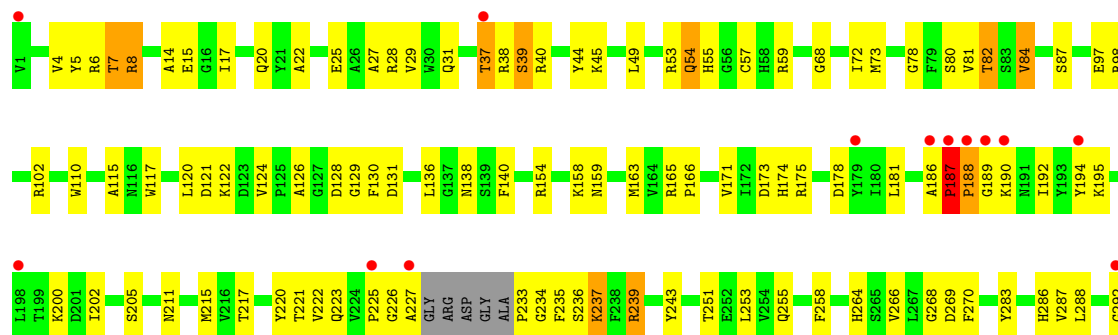


• Molecule 1: Glycine N-methyltransferase





• Molecule 1: Glycine N-methyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.43Å 84.75Å 132.25Å 90.00° 91.61° 90.00°	Depositor
Resolution (Å)	29.00 – 2.55 29.00 – 2.56	Depositor EDS
% Data completeness (in resolution range)	88.8 (29.00-2.55) 88.8 (29.00-2.56)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.57Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.240 , 0.298 0.234 , 0.290	Depositor DCC
R_{free} test set	1958 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.029 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8848	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: C2F

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	0.46	0/2189	0.91	5/2968 (0.2%)
1	B	0.49	0/2271	0.98	11/3079 (0.4%)
1	C	0.46	0/2227	0.94	10/3021 (0.3%)
1	D	0.51	0/2286	1.00	16/3100 (0.5%)
All	All	0.48	0/8973	0.96	42/12168 (0.3%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	81	VAL	N-CA-C	8.16	119.60	108.17
1	B	81	VAL	N-CA-C	8.09	119.50	108.17
1	D	202	ILE	N-CA-C	7.31	118.43	107.75
1	B	202	ILE	N-CA-C	7.06	118.50	107.28
1	A	81	VAL	N-CA-C	6.62	118.94	108.87
1	C	188	PRO	N-CA-CB	6.56	110.13	103.25
1	D	225	PRO	N-CA-CB	6.34	109.91	103.25
1	D	243	TYR	N-CA-C	-6.34	101.80	109.72
1	B	225	PRO	N-CA-CB	6.31	109.88	103.25
1	D	7	THR	N-CA-C	-6.29	104.50	111.36
1	C	81	VAL	N-CA-C	6.25	118.37	108.87
1	C	271	LYS	CA-C-N	-6.14	113.24	120.11
1	C	271	LYS	C-N-CA	-6.14	113.24	120.11
1	D	233	PRO	N-CA-CB	6.10	109.71	103.00
1	B	243	TYR	N-CA-C	-6.00	102.22	109.72
1	C	280	VAL	CA-C-N	5.98	125.94	119.78
1	C	280	VAL	C-N-CA	5.98	125.94	119.78
1	A	280	VAL	CA-C-N	5.96	125.92	119.78
1	A	280	VAL	C-N-CA	5.96	125.92	119.78
1	D	235	PHE	N-CA-C	5.96	118.62	108.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	78	GLY	N-CA-C	5.89	123.17	115.40
1	B	22	ALA	N-CA-C	5.77	119.14	111.75
1	D	22	ALA	N-CA-C	5.66	118.99	111.75
1	B	237	LYS	N-CA-C	5.57	117.26	108.96
1	B	78	GLY	N-CA-C	5.47	122.64	115.36
1	A	90	MET	N-CA-C	-5.37	104.85	111.40
1	B	178	ASP	N-CA-C	-5.36	105.52	111.36
1	D	6	ARG	N-CA-C	5.30	118.57	110.20
1	C	239	ARG	N-CA-C	5.29	117.43	109.23
1	B	37	THR	N-CA-C	-5.22	106.75	113.01
1	D	37	THR	N-CA-C	-5.17	106.80	113.01
1	D	187	PRO	CA-C-N	5.17	126.30	119.84
1	D	187	PRO	C-N-CA	5.17	126.30	119.84
1	C	200	LYS	N-CA-C	5.16	116.58	107.61
1	D	178	ASP	N-CA-C	-5.14	105.76	111.36
1	B	187	PRO	CA-C-N	5.13	126.26	119.84
1	B	187	PRO	C-N-CA	5.13	126.26	119.84
1	C	69	VAL	N-CA-C	5.13	116.47	110.62
1	C	90	MET	N-CA-C	-5.07	105.22	111.40
1	D	87	SER	N-CA-C	5.04	116.86	107.99
1	D	237	LYS	N-CA-C	5.04	116.46	108.96
1	A	24	GLY	N-CA-C	-5.02	104.96	112.89

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	2140	0	2075	89	0
1	B	2219	0	2165	83	1
1	C	2176	0	2108	91	0
1	D	2233	0	2168	73	0
2	A	33	0	23	17	0
2	B	33	0	23	11	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	5	0	0	0	0
3	C	4	0	0	0	0
3	D	4	0	0	0	0
All	All	8848	0	8562	333	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 19.

All (333) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1410:C2F:H72	1:B:207:LEU:CD2	1.97	0.95
2:A:1410:C2F:H72	1:B:207:LEU:HD22	1.50	0.90
2:A:1410:C2F:H13	1:B:239:ARG:HH12	1.42	0.84
1:D:264:HIS:HE1	1:D:286:HIS:HD2	1.26	0.83
1:B:174:HIS:CE1	1:B:286:HIS:HE1	1.97	0.82
1:A:5:TYR:CZ	2:A:1410:C2F:HB2	2.15	0.81
1:D:174:HIS:CE1	1:D:286:HIS:HE1	1.97	0.81
1:B:8:ARG:HH21	1:B:13:ALA:HA	1.45	0.81
1:D:215:MET:HE3	1:D:217:THR:OG1	1.79	0.81
1:A:274:ARG:O	1:A:277:GLN:HB2	1.81	0.81
1:B:215:MET:HE3	1:B:217:THR:OG1	1.81	0.81
1:B:264:HIS:HE1	1:B:286:HIS:HD2	1.28	0.80
1:A:220:TYR:O	1:A:237:LYS:HB2	1.81	0.80
1:C:264:HIS:CD2	1:C:265:SER:H	1.99	0.80
1:C:274:ARG:O	1:C:277:GLN:HB2	1.83	0.79
1:B:49:LEU:O	1:B:53:ARG:HG2	1.83	0.78
1:B:264:HIS:CE1	1:B:286:HIS:HD2	2.04	0.76
1:A:251:THR:HG23	1:A:264:HIS:CE1	2.21	0.76
1:D:49:LEU:O	1:D:53:ARG:HG2	1.86	0.75
1:D:264:HIS:CE1	1:D:286:HIS:HD2	2.04	0.75
1:C:55:HIS:CE1	1:C:289:LYS:HE2	2.21	0.75
1:D:264:HIS:HE1	1:D:286:HIS:CD2	2.04	0.75
1:B:264:HIS:HE1	1:B:286:HIS:CD2	2.05	0.75
1:D:174:HIS:HE1	1:D:286:HIS:HE1	1.33	0.74
1:C:120:LEU:HG	1:C:163:MET:HE3	1.70	0.74
1:B:174:HIS:HE1	1:B:286:HIS:HE1	1.34	0.74
1:B:3:SER:HB2	2:B:1420:C2F:O4	1.86	0.74
1:B:5:TYR:CD1	2:B:1420:C2F:H92	2.22	0.73
2:B:1420:C2F:H111	2:B:1420:C2F:N10	2.03	0.72
1:B:166:PRO:HB3	1:B:292:GLY:HA2	1.70	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1410:C2F:N10	2:A:1410:C2F:H111	2.05	0.71
1:B:129:GLY:HA3	1:B:165:ARG:HB3	1.72	0.70
1:A:72:ILE:CD1	1:A:97:GLU:HG2	2.22	0.70
1:A:120:LEU:HG	1:A:163:MET:HE3	1.72	0.70
1:D:68:GLY:O	1:D:72:ILE:HG13	1.92	0.69
1:C:72:ILE:CD1	1:C:97:GLU:HG2	2.22	0.69
1:A:149:ASP:OD2	1:A:151:SER:HB3	1.91	0.69
1:D:166:PRO:HB3	1:D:292:GLY:HA2	1.74	0.69
1:C:149:ASP:OD2	1:C:151:SER:HB3	1.93	0.69
1:C:48:LEU:HD23	1:C:73:MET:HE1	1.75	0.69
1:C:258:PHE:CE2	1:C:288:LEU:HD22	2.28	0.68
1:A:258:PHE:CE2	1:A:288:LEU:HD22	2.29	0.68
1:D:129:GLY:HA3	1:D:165:ARG:HB3	1.76	0.67
1:A:48:LEU:HD23	1:A:73:MET:HE1	1.76	0.67
1:C:72:ILE:HD11	1:C:97:GLU:HG2	1.77	0.67
1:A:50:GLY:O	1:A:54:GLN:HG2	1.94	0.66
1:C:50:GLY:O	1:C:54:GLN:HG2	1.95	0.66
1:C:66:GLY:HA2	1:C:90:MET:HG2	1.78	0.66
1:B:2:ASP:OD2	1:D:8:ARG:HD3	1.95	0.66
1:B:8:ARG:NH2	1:B:13:ALA:HA	2.10	0.66
1:B:68:GLY:O	1:B:72:ILE:HG13	1.96	0.65
1:A:72:ILE:HD11	1:A:97:GLU:HG2	1.79	0.65
1:D:25:GLU:HG3	1:D:28:ARG:HH21	1.62	0.64
1:A:119:THR:HB	1:A:122:LYS:HE2	1.78	0.64
2:A:1410:C2F:H13	1:B:239:ARG:NH1	2.12	0.64
1:D:220:TYR:O	1:D:237:LYS:HB2	1.98	0.64
1:B:220:TYR:O	1:B:237:LYS:HB2	1.97	0.64
1:C:119:THR:HB	1:C:122:LYS:HE2	1.79	0.64
1:A:66:GLY:HA2	1:A:90:MET:HG2	1.81	0.63
1:D:174:HIS:HE1	1:D:286:HIS:CE1	2.17	0.63
1:A:65:CYS:SG	1:A:83:SER:HB3	2.39	0.63
1:C:65:CYS:SG	1:C:83:SER:HB3	2.39	0.62
1:B:25:GLU:HG3	1:B:28:ARG:HH21	1.64	0.62
1:C:48:LEU:HD23	1:C:73:MET:CE	2.30	0.61
1:A:161:ALA:C	1:A:163:MET:H	2.09	0.61
2:B:1420:C2F:H91	1:C:207:LEU:HD22	1.83	0.61
1:B:29:VAL:HG21	1:B:236:SER:HB2	1.83	0.60
1:B:174:HIS:HE1	1:B:286:HIS:CE1	2.18	0.60
2:B:1420:C2F:H111	2:B:1420:C2F:HN1	1.66	0.60
1:C:161:ALA:C	1:C:163:MET:H	2.10	0.60
1:A:215:MET:HG2	1:A:216:VAL:N	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1420:C2F:HB2	1:D:5:TYR:CZ	2.37	0.60
1:B:4:VAL:HB	1:D:4:VAL:HB	1.82	0.59
1:D:29:VAL:HG21	1:D:236:SER:HB2	1.83	0.59
1:C:247:LEU:HD22	1:C:284:PHE:CE1	2.38	0.59
1:D:25:GLU:HG3	1:D:28:ARG:NH2	2.17	0.59
1:B:38:ARG:O	1:B:39:SER:HB2	2.03	0.58
1:A:85:ASP:HB3	1:A:91:LEU:HD13	1.85	0.58
1:A:172:ILE:HG23	1:A:172:ILE:O	2.03	0.58
1:A:48:LEU:HD23	1:A:73:MET:CE	2.32	0.58
1:A:223:GLN:HA	1:A:235:PHE:N	2.19	0.58
1:D:38:ARG:O	1:D:39:SER:HB2	2.04	0.58
2:A:1410:C2F:H111	2:A:1410:C2F:HN1	1.69	0.57
1:B:174:HIS:CE1	1:B:286:HIS:CE1	2.88	0.57
1:C:183:THR:C	1:C:185:CYS:H	2.12	0.57
1:C:247:LEU:HD22	1:C:284:PHE:CZ	2.39	0.57
1:B:174:HIS:HD2	1:B:175:ARG:O	1.87	0.57
1:D:174:HIS:HD2	1:D:175:ARG:O	1.86	0.57
2:A:1410:C2F:N10	1:B:215:MET:HE1	2.18	0.57
1:C:51:LEU:HD11	1:C:169:LEU:HD21	1.86	0.57
1:A:61:LEU:HD12	1:A:82:THR:O	2.04	0.57
1:D:174:HIS:CE1	1:D:286:HIS:CE1	2.87	0.57
1:A:287:VAL:C	1:A:288:LEU:HD12	2.29	0.57
1:A:183:THR:C	1:A:185:CYS:H	2.13	0.57
1:C:40:ARG:NH1	1:C:193:TYR:HB3	2.20	0.56
1:C:287:VAL:C	1:C:288:LEU:HD12	2.29	0.56
1:A:247:LEU:HD22	1:A:284:PHE:CE1	2.40	0.56
1:A:247:LEU:HD22	1:A:284:PHE:CZ	2.40	0.56
1:A:262:CYS:SG	1:A:288:LEU:HD23	2.45	0.56
1:A:5:TYR:OH	2:A:1410:C2F:HB2	2.05	0.56
1:B:25:GLU:HG3	1:B:28:ARG:NH2	2.19	0.56
1:C:6:ARG:O	1:D:239:ARG:HD3	2.05	0.56
1:A:264:HIS:CD2	1:A:265:SER:N	2.74	0.56
2:A:1410:C2F:H71	1:C:5:TYR:CD1	2.40	0.56
1:B:54:GLN:O	1:B:54:GLN:HG2	2.04	0.56
1:C:172:ILE:HG23	1:C:172:ILE:O	2.05	0.55
1:B:136:LEU:HA	1:B:173:ASP:OD1	2.06	0.55
1:C:61:LEU:HD12	1:C:82:THR:O	2.06	0.55
1:C:258:PHE:HE2	1:C:288:LEU:HD22	1.71	0.55
1:A:221:THR:HG22	1:A:222:VAL:N	2.21	0.55
1:A:192:ILE:HG21	1:A:283:TYR:CE1	2.42	0.55
1:A:241:SER:HB2	1:B:7:THR:HG23	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:LYS:HG2	1:B:73:MET:HE2	1.89	0.54
2:B:1420:C2F:O	2:B:1420:C2F:HG1	2.08	0.54
1:C:85:ASP:HB3	1:C:91:LEU:HD13	1.89	0.54
1:D:45:LYS:HG2	1:D:73:MET:HE2	1.90	0.54
1:C:215:MET:HG2	1:C:216:VAL:N	2.22	0.54
1:A:177:TYR:HB2	1:A:245:HIS:O	2.08	0.54
1:A:72:ILE:HD13	1:A:97:GLU:HG2	1.89	0.53
1:D:258:PHE:CE2	1:D:288:LEU:HD22	2.44	0.53
1:A:156:ALA:O	1:A:160:ILE:HG13	2.09	0.53
1:C:215:MET:HE3	1:C:217:THR:OG1	2.08	0.53
1:A:135:CYS:HB3	1:A:172:ILE:HG13	1.91	0.53
1:D:54:GLN:O	1:D:54:GLN:HG2	2.08	0.53
2:A:1410:C2F:H92	1:C:5:TYR:CD1	2.44	0.52
1:A:51:LEU:HD11	1:A:169:LEU:HD21	1.89	0.52
1:A:251:THR:HG23	1:A:264:HIS:HE1	1.71	0.52
1:C:268:GLY:O	1:C:269:ASP:HB2	2.09	0.52
1:A:258:PHE:HE2	1:A:288:LEU:HD22	1.73	0.52
2:A:1410:C2F:O	2:A:1410:C2F:HG1	2.10	0.52
1:B:45:LYS:HG2	1:B:73:MET:CE	2.39	0.52
2:B:1420:C2F:H13	1:C:239:ARG:HH12	1.74	0.52
1:B:84:VAL:CG1	1:B:115:ALA:HB3	2.40	0.51
1:D:84:VAL:HG13	1:D:115:ALA:HB3	1.91	0.51
2:A:1410:C2F:C13	1:B:239:ARG:HH12	2.20	0.51
1:C:72:ILE:HD13	1:C:97:GLU:HG2	1.90	0.51
1:A:221:THR:CG2	1:A:222:VAL:N	2.73	0.51
2:B:1420:C2F:C14	1:C:215:MET:HE1	2.41	0.51
1:C:177:TYR:HB2	1:C:245:HIS:O	2.09	0.51
1:A:6:ARG:O	1:B:239:ARG:HD3	2.11	0.51
1:A:268:GLY:O	1:A:269:ASP:HB2	2.11	0.51
1:C:136:LEU:HA	1:C:173:ASP:OD1	2.11	0.51
1:C:173:ASP:OD2	1:C:193:TYR:OH	2.17	0.51
1:B:14:ALA:HB3	1:B:17:ILE:HD11	1.93	0.51
1:A:136:LEU:HA	1:A:173:ASP:OD1	2.10	0.51
1:B:6:ARG:NH1	1:B:10:LEU:HD21	2.26	0.51
1:A:215:MET:HE3	1:A:217:THR:OG1	2.11	0.50
1:A:149:ASP:C	1:A:151:SER:H	2.19	0.50
1:B:84:VAL:HG13	1:B:115:ALA:HB3	1.94	0.50
1:D:120:LEU:HB3	1:D:159:ASN:HB3	1.92	0.50
1:D:258:PHE:HE2	1:D:288:LEU:HD22	1.75	0.50
1:C:135:CYS:HB3	1:C:172:ILE:HG13	1.93	0.50
1:C:250:PHE:HD2	1:C:286:HIS:CE1	2.30	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:ALA:O	1:D:128:ASP:OD1	2.29	0.50
1:A:75:VAL:HG23	1:A:81:VAL:HG21	1.94	0.50
1:B:37:THR:HG23	1:B:194:TYR:CD2	2.47	0.49
1:B:59:ARG:HA	1:B:80:SER:HB3	1.94	0.49
1:C:264:HIS:CD2	1:C:265:SER:N	2.77	0.49
1:C:222:VAL:HG22	1:C:223:GLN:N	2.27	0.49
1:A:207:LEU:CD2	2:B:1420:C2F:HA	2.42	0.49
1:C:149:ASP:C	1:C:151:SER:H	2.20	0.49
1:B:120:LEU:HB3	1:B:159:ASN:HB3	1.94	0.49
1:B:121:ASP:OD1	1:B:122:LYS:HG3	2.13	0.49
1:C:51:LEU:O	1:C:55:HIS:HD2	1.95	0.49
1:D:136:LEU:HA	1:D:173:ASP:OD1	2.13	0.49
1:C:156:ALA:O	1:C:160:ILE:HG13	2.12	0.49
1:D:223:GLN:HA	1:D:234:GLY:O	2.13	0.49
1:A:37:THR:O	1:A:37:THR:HG23	2.11	0.49
1:A:59:ARG:HA	1:A:80:SER:HB3	1.94	0.49
1:D:126:ALA:C	1:D:128:ASP:H	2.20	0.49
1:B:72:ILE:HD13	1:B:97:GLU:HG2	1.95	0.49
1:D:37:THR:HG23	1:D:194:TYR:CD2	2.47	0.48
1:D:45:LYS:HG2	1:D:73:MET:CE	2.42	0.48
1:A:179:TYR:HE1	1:A:187:PRO:HD3	1.78	0.48
1:B:136:LEU:HD21	1:B:171:VAL:HG12	1.95	0.48
1:D:84:VAL:CG1	1:D:115:ALA:HB3	2.42	0.48
1:A:41:THR:HG22	1:A:43:GLU:H	1.77	0.48
1:C:37:THR:O	1:C:37:THR:HG23	2.12	0.48
1:C:269:ASP:O	1:C:270:PHE:HB2	2.13	0.48
1:B:188:PRO:O	1:B:190:LYS:N	2.46	0.48
1:B:258:PHE:CE2	1:B:288:LEU:HD22	2.47	0.48
1:B:258:PHE:HE2	1:B:288:LEU:HD22	1.79	0.48
1:C:59:ARG:HA	1:C:80:SER:HB3	1.95	0.48
1:A:154:ARG:HG2	1:A:154:ARG:HH11	1.79	0.48
1:D:14:ALA:HB3	1:D:17:ILE:HD11	1.95	0.48
1:D:39:SER:OG	1:D:195:LYS:HE3	2.14	0.47
1:A:250:PHE:HD2	1:A:286:HIS:CE1	2.32	0.47
1:B:192:ILE:HD13	1:B:270:PHE:CZ	2.49	0.47
1:C:223:GLN:HA	1:C:235:PHE:N	2.29	0.47
1:C:41:THR:HG22	1:C:43:GLU:H	1.79	0.47
1:D:53:ARG:O	1:D:55:HIS:N	2.47	0.47
1:A:150:GLN:NE2	1:A:243:TYR:OH	2.47	0.47
1:A:174:HIS:ND1	1:A:174:HIS:O	2.48	0.47
1:D:192:ILE:HD13	1:D:270:PHE:CZ	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:ARG:C	1:B:55:HIS:H	2.23	0.47
1:D:136:LEU:HD21	1:D:171:VAL:HG12	1.97	0.47
1:B:83:SER:O	1:B:84:VAL:HG23	2.15	0.47
1:B:188:PRO:C	1:B:190:LYS:H	2.23	0.47
1:C:55:HIS:ND1	1:C:289:LYS:HE2	2.30	0.47
2:A:1410:C2F:HN1	1:B:215:MET:HE1	1.81	0.47
1:D:72:ILE:HD13	1:D:97:GLU:HG2	1.95	0.47
1:A:192:ILE:HD13	1:A:283:TYR:CD1	2.50	0.46
1:D:195:LYS:H	1:D:195:LYS:HG2	1.55	0.46
1:B:40:ARG:HB3	1:B:44:TYR:HD2	1.81	0.46
1:D:188:PRO:C	1:D:190:LYS:H	2.23	0.46
1:A:264:HIS:HD2	1:A:265:SER:H	1.64	0.46
1:C:154:ARG:HH11	1:C:154:ARG:HG2	1.81	0.46
1:C:51:LEU:HD12	1:C:55:HIS:CD2	2.51	0.46
1:A:136:LEU:HD21	1:A:171:VAL:HG12	1.98	0.46
1:B:53:ARG:O	1:B:55:HIS:N	2.49	0.46
1:C:75:VAL:HG23	1:C:81:VAL:HG21	1.98	0.46
1:D:188:PRO:O	1:D:190:LYS:N	2.48	0.46
1:B:124:VAL:HG21	1:B:163:MET:HE2	1.97	0.46
1:A:220:TYR:O	1:A:237:LYS:CB	2.61	0.45
1:A:203:THR:HG23	1:C:210:ASN:HD21	1.80	0.45
1:B:126:ALA:C	1:B:128:ASP:H	2.23	0.45
1:D:53:ARG:C	1:D:55:HIS:H	2.24	0.45
1:D:121:ASP:OD1	1:D:122:LYS:HG3	2.17	0.45
1:C:176:ASN:O	1:C:180:ILE:HG13	2.17	0.45
1:A:161:ALA:C	1:A:163:MET:N	2.75	0.45
1:A:264:HIS:CD2	1:A:265:SER:H	2.34	0.45
1:B:82:THR:CG2	1:B:130:PHE:HE2	2.30	0.45
1:B:82:THR:HG21	1:B:130:PHE:HE2	1.82	0.45
1:A:172:ILE:O	1:A:172:ILE:CG2	2.65	0.45
2:B:1420:C2F:HB2	1:D:5:TYR:CE1	2.52	0.45
1:D:221:THR:HA	1:D:237:LYS:HB3	1.99	0.45
1:B:221:THR:HA	1:B:237:LYS:HB3	2.00	0.44
1:C:269:ASP:O	1:C:270:PHE:CB	2.66	0.44
1:D:40:ARG:HB3	1:D:44:TYR:HD2	1.81	0.44
1:B:39:SER:OG	1:B:195:LYS:HE3	2.18	0.44
1:B:268:GLY:O	1:B:269:ASP:HB2	2.17	0.44
1:C:183:THR:O	1:C:185:CYS:N	2.50	0.44
1:A:176:ASN:O	1:A:180:ILE:HG13	2.17	0.44
1:D:175:ARG:HG2	1:D:283:TYR:CE2	2.53	0.44
1:D:268:GLY:O	1:D:269:ASP:HB2	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:ALA:HB1	1:B:258:PHE:CE1	2.53	0.44
1:C:35:GLY:O	1:C:38:ARG:HG2	2.18	0.44
1:D:200:LYS:HG2	1:D:222:VAL:HG22	2.00	0.44
1:B:195:LYS:H	1:B:195:LYS:HG2	1.54	0.44
1:C:262:CYS:SG	1:C:288:LEU:HD23	2.58	0.44
1:A:203:THR:HG23	1:C:210:ASN:ND2	2.33	0.43
1:C:172:ILE:O	1:C:172:ILE:CG2	2.65	0.43
1:C:209:VAL:O	1:C:210:ASN:HB2	2.18	0.43
1:C:268:GLY:N	1:C:271:LYS:O	2.44	0.43
1:C:40:ARG:HD3	1:C:40:ARG:HA	1.83	0.43
1:C:174:HIS:O	1:C:174:HIS:ND1	2.51	0.43
1:A:33:TYR:O	1:A:37:THR:HG22	2.19	0.43
1:D:251:THR:HG22	1:D:255:GLN:HE21	1.83	0.43
1:B:226:GLY:O	1:B:227:ALA:C	2.61	0.43
1:D:226:GLY:O	1:D:227:ALA:C	2.62	0.43
1:A:29:VAL:HG21	1:A:236:SER:HB3	2.01	0.43
1:A:138:ASN:CG	1:A:175:ARG:HG3	2.43	0.43
1:B:98:ARG:HB2	1:B:110:TRP:CE3	2.53	0.43
1:C:188:PRO:HA	1:C:202:ILE:HD12	2.01	0.43
1:D:124:VAL:HG21	1:D:163:MET:HE2	2.00	0.43
1:A:61:LEU:HB2	1:A:130:PHE:CD2	2.54	0.43
1:C:136:LEU:HD21	1:C:171:VAL:HG12	2.01	0.43
1:C:150:GLN:NE2	1:C:243:TYR:OH	2.51	0.43
1:B:200:LYS:HG2	1:B:222:VAL:HG22	2.01	0.43
1:C:187:PRO:HB2	1:C:190:LYS:CB	2.48	0.43
1:D:82:THR:CG2	1:D:130:PHE:HE2	2.32	0.43
1:A:212:LYS:O	1:A:213:ALA:C	2.62	0.42
1:A:209:VAL:O	1:A:210:ASN:HB2	2.19	0.42
1:C:220:TYR:O	1:C:237:LYS:HB2	2.18	0.42
1:A:195:LYS:O	1:A:195:LYS:HG3	2.20	0.42
1:B:251:THR:O	1:B:255:GLN:HG3	2.19	0.42
1:C:116:ASN:O	1:C:117:TRP:C	2.61	0.42
1:D:117:TRP:CH2	1:D:140:PHE:HA	2.55	0.42
1:A:265:SER:HB3	1:A:287:VAL:HG22	2.01	0.42
1:B:175:ARG:HG2	1:B:283:TYR:CE2	2.55	0.42
1:C:33:TYR:O	1:C:37:THR:HG22	2.19	0.42
1:D:98:ARG:HB2	1:D:110:TRP:CE3	2.54	0.42
1:A:92:LYS:HD2	1:B:114:GLU:OE1	2.18	0.42
1:A:269:ASP:O	1:A:270:PHE:HB2	2.19	0.42
1:D:40:ARG:HH11	1:D:194:TYR:HB3	1.85	0.42
1:A:5:TYR:HB3	2:A:1410:C2F:C13	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:149:ASP:C	1:C:151:SER:N	2.78	0.42
1:A:183:THR:O	1:A:185:CYS:N	2.51	0.42
1:C:74:LEU:O	1:C:79:PHE:HB2	2.20	0.42
1:D:82:THR:HG21	1:D:130:PHE:HE2	1.84	0.42
1:C:137:GLY:O	1:C:138:ASN:HB3	2.20	0.42
1:A:183:THR:C	1:A:185:CYS:N	2.77	0.42
1:A:41:THR:HG22	1:A:43:GLU:N	2.34	0.41
1:A:137:GLY:O	1:A:138:ASN:HB3	2.19	0.41
1:B:208:THR:O	1:D:205:SER:HA	2.20	0.41
1:B:53:ARG:C	1:B:55:HIS:N	2.78	0.41
1:B:251:THR:HG22	1:B:255:GLN:HE21	1.84	0.41
1:D:59:ARG:HA	1:D:80:SER:HB3	2.02	0.41
1:A:5:TYR:CE2	2:A:1410:C2F:HB2	2.53	0.41
1:B:154:ARG:NH1	1:B:253:LEU:HD11	2.34	0.41
1:C:161:ALA:C	1:C:163:MET:N	2.76	0.41
1:A:149:ASP:C	1:A:151:SER:N	2.77	0.41
2:A:1410:C2F:H71	1:C:5:TYR:HD1	1.82	0.41
1:A:116:ASN:O	1:A:117:TRP:C	2.63	0.41
1:A:221:THR:O	1:A:222:VAL:HG23	2.21	0.41
1:C:43:GLU:H	1:C:43:GLU:HG3	1.71	0.41
1:D:53:ARG:C	1:D:55:HIS:N	2.78	0.41
1:D:138:ASN:CG	1:D:175:ARG:HG3	2.46	0.41
1:D:186:ALA:O	1:D:187:PRO:O	2.39	0.41
1:D:251:THR:O	1:D:255:GLN:HG3	2.20	0.41
1:A:65:CYS:HB3	1:A:85:ASP:HB2	2.02	0.41
1:B:211:ASN:HD22	1:B:211:ASN:HA	1.71	0.41
1:C:168:GLY:O	1:C:289:LYS:HA	2.21	0.41
1:C:251:THR:O	1:C:255:GLN:HG3	2.21	0.41
1:B:98:ARG:HB2	1:B:110:TRP:CZ3	2.56	0.41
1:D:138:ASN:ND2	1:D:175:ARG:HG3	2.36	0.41
1:D:57:CYS:HA	1:D:131:ASP:HB3	2.03	0.41
1:A:91:LEU:O	1:A:92:LYS:C	2.64	0.40
1:A:130:PHE:O	1:A:165:ARG:HG2	2.21	0.40
1:A:187:PRO:HA	1:A:188:PRO:HD2	1.96	0.40
1:B:280:VAL:HA	1:B:281:PRO:HD3	1.87	0.40
1:C:41:THR:HG22	1:C:43:GLU:N	2.35	0.40
1:C:277:GLN:HG2	1:C:278:ALA:H	1.86	0.40
1:B:40:ARG:HH11	1:B:194:TYR:HB3	1.86	0.40
1:C:55:HIS:HE1	1:C:289:LYS:HE2	1.80	0.40
1:D:264:HIS:CE1	1:D:286:HIS:CD2	2.90	0.40
1:A:40:ARG:HA	1:A:40:ARG:HD3	1.81	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:ALA:O	1:B:187:PRO:O	2.39	0.40
1:C:65:CYS:HB3	1:C:85:ASP:HB2	2.03	0.40
1:C:84:VAL:CG1	1:C:115:ALA:HB3	2.51	0.40
1:D:29:VAL:HG21	1:D:236:SER:CB	2.51	0.40
1:D:154:ARG:NH1	1:D:253:LEU:HD11	2.36	0.40
1:D:211:ASN:HD22	1:D:211:ASN:HA	1.72	0.40
1:B:138:ASN:O	1:B:138:ASN:CG	2.65	0.40
1:C:130:PHE:O	1:C:165:ARG:HG2	2.21	0.40
1:C:138:ASN:CG	1:C:175:ARG:HG3	2.47	0.40
1:D:27:ALA:O	1:D:31:GLN:HG3	2.21	0.40
1:A:143:LEU:HD12	1:A:144:PRO:HD2	2.04	0.40
1:B:72:ILE:CD1	1:B:97:GLU:HG2	2.51	0.40
1:B:117:TRP:CH2	1:B:140:PHE:HA	2.57	0.40
1:C:9:SER:HB2	1:C:12:VAL:CG2	2.51	0.40
1:C:9:SER:HB2	1:C:12:VAL:HG21	2.03	0.40
1:C:104:GLU:O	1:C:105:PRO:C	2.65	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:102:ARG:NH2	1:B:276:GLY:O[2_555]	2.08	0.12

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	271/292 (93%)	233 (86%)	31 (11%)	7 (3%)	4	3
1	B	281/292 (96%)	264 (94%)	11 (4%)	6 (2%)	5	5
1	C	277/292 (95%)	239 (86%)	33 (12%)	5 (2%)	6	7
1	D	283/292 (97%)	265 (94%)	12 (4%)	6 (2%)	5	5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1112/1168 (95%)	1001 (90%)	87 (8%)	24 (2%)	5 5

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	291	THR
1	B	20	GLN
1	C	291	THR
1	D	20	GLN
1	A	195	LYS
1	B	39	SER
1	B	54	GLN
1	B	187	PRO
1	B	189	GLY
1	D	39	SER
1	D	54	GLN
1	D	187	PRO
1	D	189	GLY
1	A	191	ASN
1	A	162	SER
1	B	188	PRO
1	C	192	ILE
1	D	188	PRO
1	A	55	HIS
1	C	162	SER
1	C	55	HIS
1	C	184	GLY
1	A	184	GLY
1	A	268	GLY

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	222/242 (92%)	216 (97%)	6 (3%)	39 58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	233/242 (96%)	221 (95%)	12 (5%)	21	32
1	C	225/242 (93%)	219 (97%)	6 (3%)	39	58
1	D	234/242 (97%)	222 (95%)	12 (5%)	21	33
All	All	914/968 (94%)	878 (96%)	36 (4%)	28	43

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	84	VAL
1	A	91	LEU
1	A	92	LYS
1	A	102	ARG
1	A	208	THR
1	B	7	THR
1	B	15	GLU
1	B	82	THR
1	B	84	VAL
1	B	102	ARG
1	B	158	LYS
1	B	181	LEU
1	B	187	PRO
1	B	211	ASN
1	B	239	ARG
1	B	266	VAL
1	B	287	VAL
1	C	31	GLN
1	C	84	VAL
1	C	91	LEU
1	C	92	LYS
1	C	102	ARG
1	C	208	THR
1	D	7	THR
1	D	8	ARG
1	D	15	GLU
1	D	82	THR
1	D	84	VAL
1	D	102	ARG
1	D	158	LYS
1	D	181	LEU
1	D	187	PRO

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Mol	Chain	Res	Type
1	D	239	ARG
1	D	266	VAL
1	D	287	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	150	GLN
1	A	245	HIS
1	A	255	GLN
1	A	286	HIS
1	B	31	GLN
1	B	54	GLN
1	B	150	GLN
1	B	174	HIS
1	B	211	ASN
1	B	255	GLN
1	B	263	GLN
1	B	264	HIS
1	B	286	HIS
1	C	31	GLN
1	C	55	HIS
1	C	150	GLN
1	C	255	GLN
1	C	264	HIS
1	C	286	HIS
1	D	31	GLN
1	D	54	GLN
1	D	174	HIS
1	D	211	ASN
1	D	255	GLN
1	D	263	GLN
1	D	264	HIS
1	D	286	HIS

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	C2F	B	1420	-	32,35,35	1.68	7 (21%)	36,49,49	2.02	7 (19%)
2	C2F	A	1410	-	32,35,35	1.60	6 (18%)	36,49,49	2.01	7 (19%)

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	C2F	B	1420	-	-	1/22/35/35	0/3/3/3
2	C2F	A	1410	-	-	1/22/35/35	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1420	C2F	C6-N5	5.20	1.54	1.48
2	A	1410	C2F	C6-N5	5.12	1.54	1.48
2	B	1420	C2F	C7-C6	3.12	1.55	1.52
2	A	1410	C2F	CA-N	2.61	1.51	1.45
2	B	1420	C2F	C17-C12	2.59	1.43	1.39
2	A	1410	C2F	C17-C12	2.54	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1420	C2F	C14-C15	2.50	1.43	1.39
2	B	1420	C2F	C13-C12	2.50	1.43	1.39
2	A	1410	C2F	C7-C6	2.46	1.54	1.52
2	B	1420	C2F	CA-N	2.35	1.50	1.45
2	A	1410	C2F	C13-C12	2.15	1.42	1.39
2	B	1420	C2F	C16-C15	2.14	1.42	1.39
2	A	1410	C2F	C14-C15	2.04	1.42	1.39

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1410	C2F	CA-N-C	5.94	135.81	121.56
2	B	1420	C2F	CA-N-C	5.85	135.60	121.56
2	B	1420	C2F	C4A-C4-N3	4.70	119.22	110.94
2	A	1410	C2F	C4A-C4-N3	4.64	119.11	110.94
2	B	1420	C2F	C2-N1-C8A	4.56	121.40	113.36
2	B	1420	C2F	CB-CA-N	4.53	119.87	110.91
2	A	1410	C2F	C2-N1-C8A	4.43	121.17	113.36
2	A	1410	C2F	CB-CA-N	4.42	119.67	110.91
2	A	1410	C2F	CT-CA-N	-3.44	102.60	110.57
2	B	1420	C2F	CT-CA-N	-3.31	102.89	110.57
2	A	1410	C2F	O4-C4-C4A	-2.49	121.50	127.62
2	B	1420	C2F	O4-C4-C4A	-2.46	121.57	127.62
2	B	1420	C2F	C2-N3-C4	-2.17	121.17	125.11
2	A	1410	C2F	C2-N3-C4	-2.03	121.42	125.11

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1420	C2F	CB-CA-N-C
2	A	1410	C2F	CB-CA-N-C

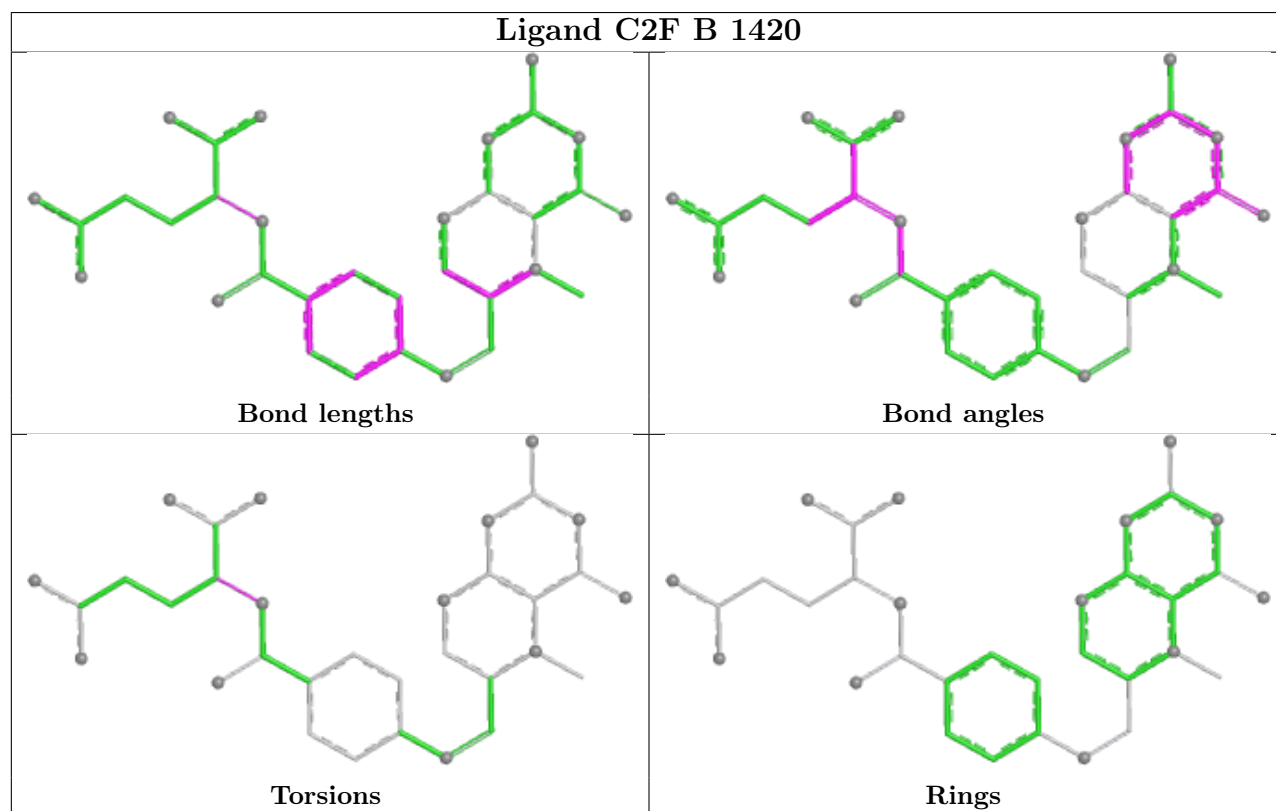
There are no ring outliers.

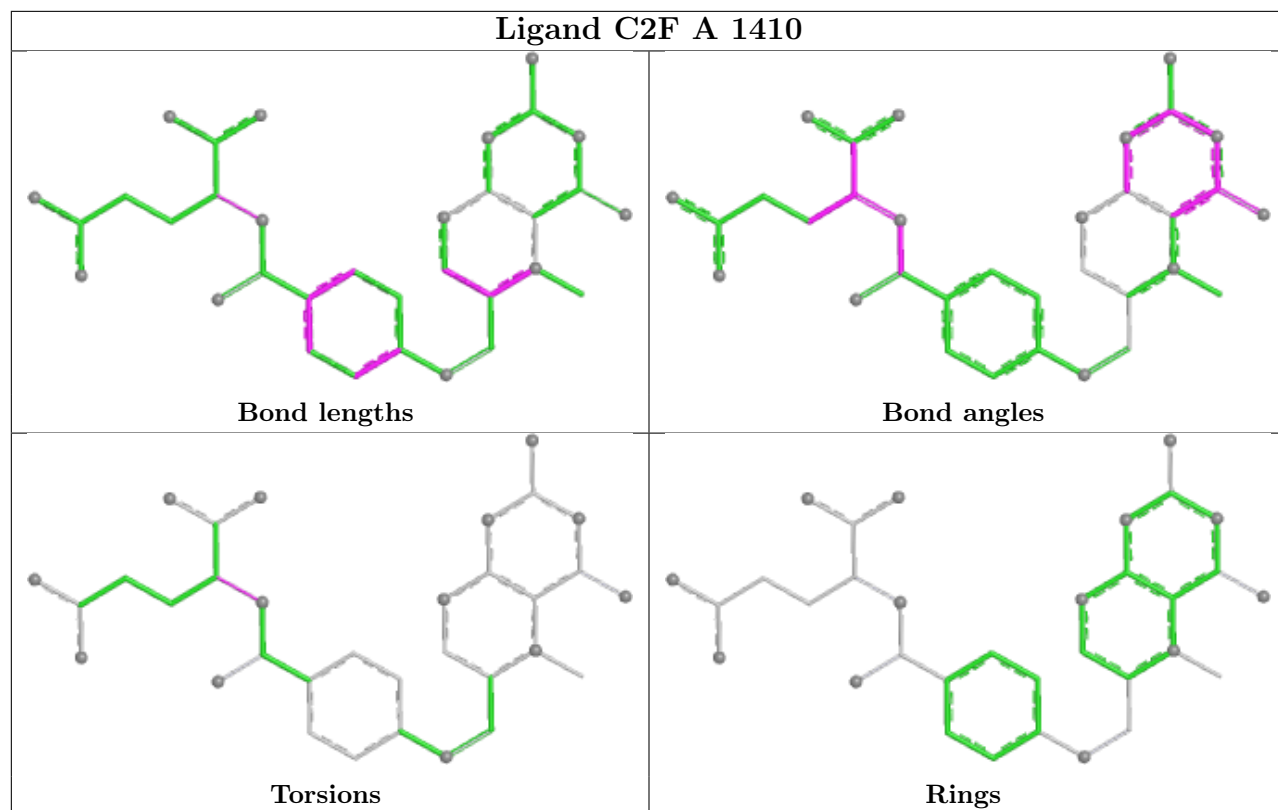
2 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1420	C2F	11	0
2	A	1410	C2F	17	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

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	277/292 (94%)	0.68	16 (5%)	29 28	22, 67, 111, 132	0
1	B	285/292 (97%)	0.25	15 (5%)	32 31	24, 48, 100, 136	0
1	C	281/292 (96%)	0.47	16 (5%)	29 29	19, 63, 110, 132	0
1	D	287/292 (98%)	0.16	13 (4%)	38 38	19, 46, 102, 136	0
All	All	1130/1168 (96%)	0.39	60 (5%)	32 31	19, 56, 108, 136	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	189	GLY	4.6
1	D	292	GLY	4.4
1	B	235	PHE	4.3
1	D	37	THR	3.9
1	B	227	ALA	3.7
1	C	187	PRO	3.5
1	A	191	ASN	3.4
1	A	51	LEU	3.3
1	A	292	GLY	3.2
1	C	58	HIS	3.1
1	A	58	HIS	3.1
1	D	186	ALA	3.1
1	B	187	PRO	3.0
1	B	224	VAL	2.9
1	A	185	CYS	2.9
1	B	190	LYS	2.9
1	C	235	PHE	2.8
1	D	188	PRO	2.8
1	B	37	THR	2.8
1	C	42	ALA	2.7
1	D	225	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	1	VAL	2.7
1	A	60	VAL	2.7
1	A	291	THR	2.7
1	D	190	LYS	2.6
1	B	1	VAL	2.5
1	C	292	GLY	2.5
1	C	37	THR	2.5
1	C	291	THR	2.5
1	A	29	VAL	2.4
1	B	223	GLN	2.4
1	A	270	PHE	2.4
1	D	189	GLY	2.4
1	C	185	CYS	2.3
1	C	57	CYS	2.3
1	D	227	ALA	2.3
1	D	194	TYR	2.3
1	A	187	PRO	2.3
1	B	188	PRO	2.3
1	C	189	GLY	2.2
1	D	198	LEU	2.2
1	B	186	ALA	2.2
1	B	5	TYR	2.2
1	C	184	GLY	2.2
1	D	187	PRO	2.2
1	A	56	GLY	2.1
1	D	179	TYR	2.1
1	A	272	PRO	2.1
1	A	55	HIS	2.1
1	A	202	ILE	2.1
1	C	128	ASP	2.1
1	C	169	LEU	2.1
1	A	183	THR	2.1
1	A	223	GLN	2.1
1	B	126	ALA	2.1
1	C	254	VAL	2.0
1	B	78	GLY	2.0
1	B	184	GLY	2.0
1	C	183	THR	2.0
1	C	202	ILE	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

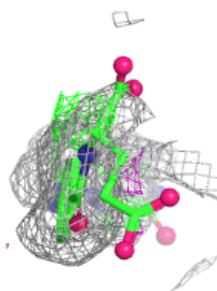
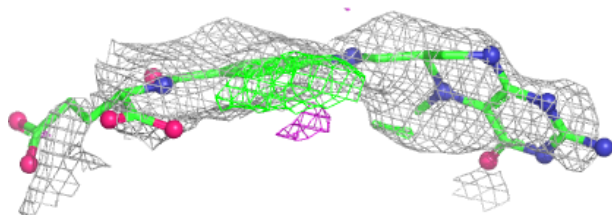
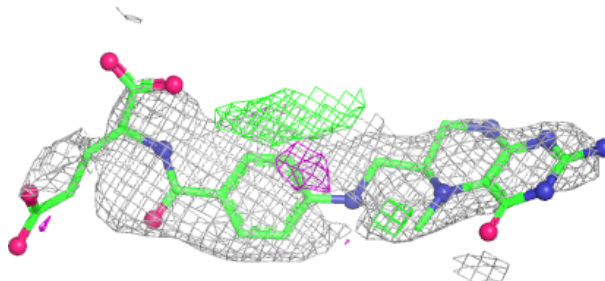
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	C2F	B	1420	33/33	0.69	0.18	68,96,112,113	0
2	C2F	A	1410	33/33	0.72	0.22	67,91,105,113	0

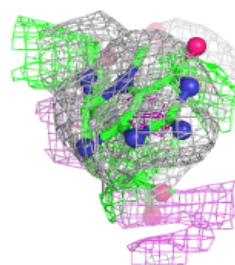
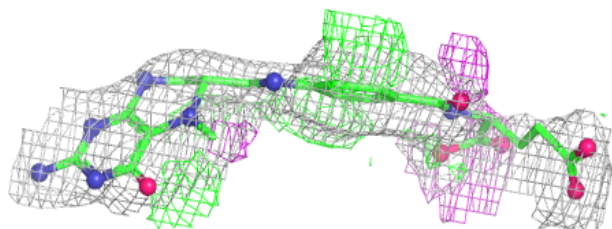
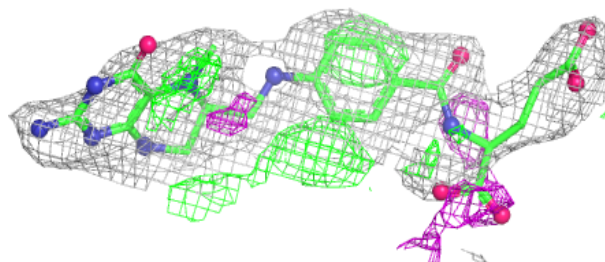
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around C2F B 1420:

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

**Electron density around C2F A 1410:**

$2mF_o-DF_c$ (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.