



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2024 – 05:37 PM EDT

PDB ID : 2P83
Title : Potent and selective isophthalamide S2 hydroxyethylamine inhibitor of BACE1
Authors : Benson, T.E.; Prince, D.B.; Tomasselli, A.G.; Emmons, T.L.; Paddock, D.J.
Deposited on : 2007-03-21
Resolution : 2.50 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 : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 2.37.1
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) : 2.37.1

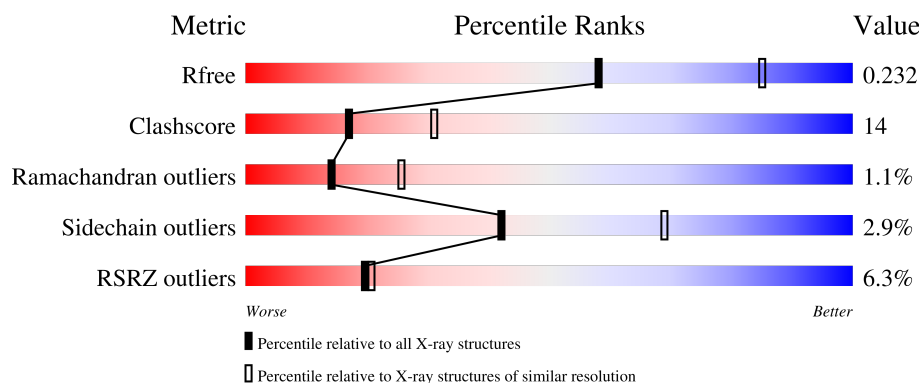
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	455	<div> <div>3%</div> <div>58%</div> <div>22%</div> <div>18%</div> </div>
1	B	455	<div> <div>6%</div> <div>59%</div> <div>21%</div> <div>18%</div> </div>
1	C	455	<div> <div>6%</div> <div>58%</div> <div>22%</div> <div>18%</div> </div>

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	PO4	B	704	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9104 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 Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	371	Total	C	N	O	S	0	0	0
			2924	1874	487	549	14			
1	B	372	Total	C	N	O	S	0	0	0
			2929	1879	487	549	14			
1	C	373	Total	C	N	O	S	0	0	0
			2937	1883	488	552	14			

There are 45 discrepancies between the modelled and reference sequences:

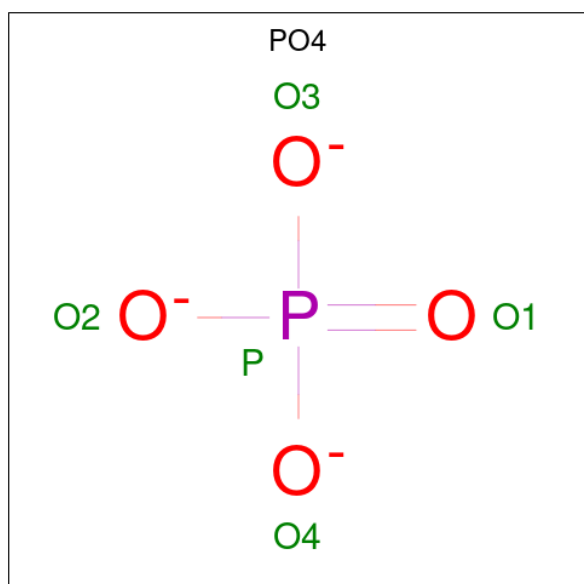
Chain	Residue	Modelled	Actual	Comment	Reference
A	1P	MET	-	EXPRESSION TAG	UNP P56817
A	2P	ALA	-	EXPRESSION TAG	UNP P56817
A	3P	SER	-	EXPRESSION TAG	UNP P56817
A	4P	MET	-	EXPRESSION TAG	UNP P56817
A	5P	THR	-	EXPRESSION TAG	UNP P56817
A	6P	GLY	-	EXPRESSION TAG	UNP P56817
A	7P	GLY	-	EXPRESSION TAG	UNP P56817
A	8P	GLN	-	EXPRESSION TAG	UNP P56817
A	9P	GLN	-	EXPRESSION TAG	UNP P56817
A	10P	MET	-	EXPRESSION TAG	UNP P56817
A	11P	GLY	-	EXPRESSION TAG	UNP P56817
A	12P	ARG	-	EXPRESSION TAG	UNP P56817
A	13P	GLY	-	EXPRESSION TAG	UNP P56817
A	14P	SER	-	EXPRESSION TAG	UNP P56817
A	15P	MET	-	EXPRESSION TAG	UNP P56817
B	1P	MET	-	EXPRESSION TAG	UNP P56817
B	2P	ALA	-	EXPRESSION TAG	UNP P56817
B	3P	SER	-	EXPRESSION TAG	UNP P56817
B	4P	MET	-	EXPRESSION TAG	UNP P56817
B	5P	THR	-	EXPRESSION TAG	UNP P56817
B	6P	GLY	-	EXPRESSION TAG	UNP P56817
B	7P	GLY	-	EXPRESSION TAG	UNP P56817
B	8P	GLN	-	EXPRESSION TAG	UNP P56817

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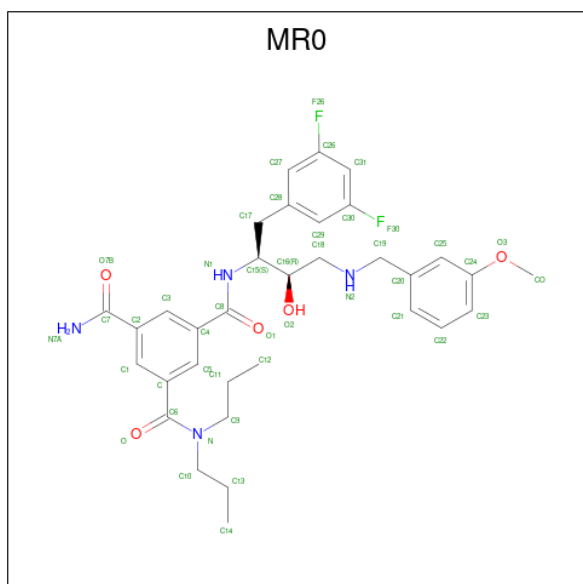
Chain	Residue	Modelled	Actual	Comment	Reference
B	9P	GLN	-	EXPRESSION TAG	UNP P56817
B	10P	MET	-	EXPRESSION TAG	UNP P56817
B	11P	GLY	-	EXPRESSION TAG	UNP P56817
B	12P	ARG	-	EXPRESSION TAG	UNP P56817
B	13P	GLY	-	EXPRESSION TAG	UNP P56817
B	14P	SER	-	EXPRESSION TAG	UNP P56817
B	15P	MET	-	EXPRESSION TAG	UNP P56817
C	1P	MET	-	EXPRESSION TAG	UNP P56817
C	2P	ALA	-	EXPRESSION TAG	UNP P56817
C	3P	SER	-	EXPRESSION TAG	UNP P56817
C	4P	MET	-	EXPRESSION TAG	UNP P56817
C	5P	THR	-	EXPRESSION TAG	UNP P56817
C	6P	GLY	-	EXPRESSION TAG	UNP P56817
C	7P	GLY	-	EXPRESSION TAG	UNP P56817
C	8P	GLN	-	EXPRESSION TAG	UNP P56817
C	9P	GLN	-	EXPRESSION TAG	UNP P56817
C	10P	MET	-	EXPRESSION TAG	UNP P56817
C	11P	GLY	-	EXPRESSION TAG	UNP P56817
C	12P	ARG	-	EXPRESSION TAG	UNP P56817
C	13P	GLY	-	EXPRESSION TAG	UNP P56817
C	14P	SER	-	EXPRESSION TAG	UNP P56817
C	15P	MET	-	EXPRESSION TAG	UNP P56817

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0

- Molecule 3 is N 3 -{(1S,2R)-1-(3,5-DIFLUOROBENZYL)-2-HYDROXY-3-[(3-METHOXY BENZYL)AMINO]PROPYL}-N 1 ,N 1 -DIPROPYLBENZENE-1,3,5-TRICARBOXAMID E (three-letter code: MR0) (formula: C₃₃H₄₀F₂N₄O₅).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C F N O 44 33 2 4 5	0	0
3	B	1	Total C F N O 44 33 2 4 5	0	0
3	C	1	Total C F N O 44 33 2 4 5	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	64	Total O 64 64	0	0
4	B	52	Total O 52 52	0	0

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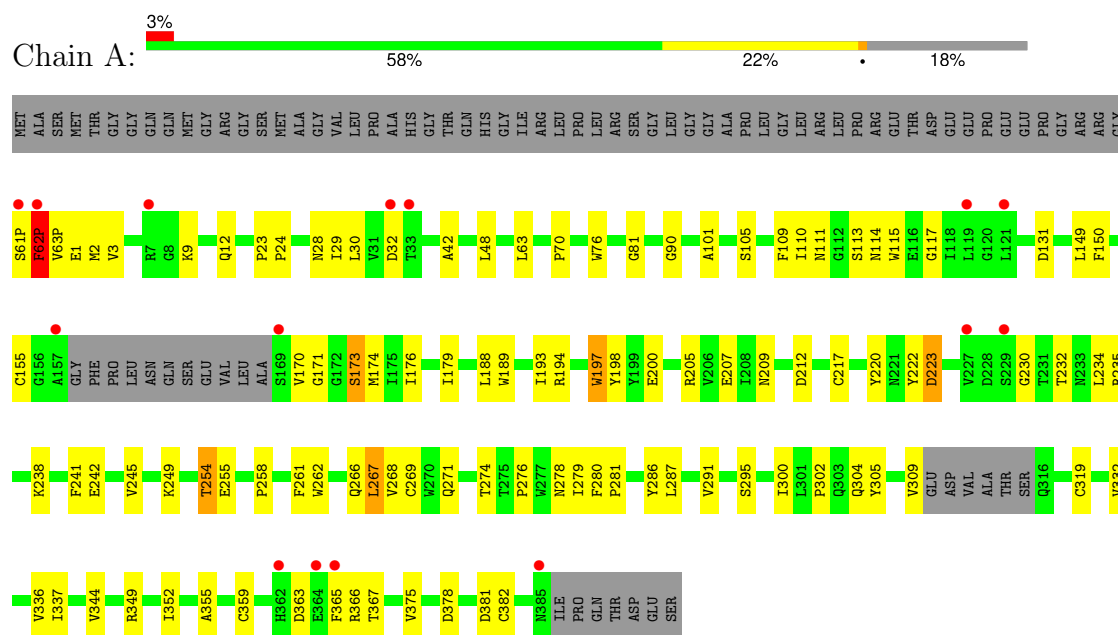
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	46	Total	O	0	0
			46	46		

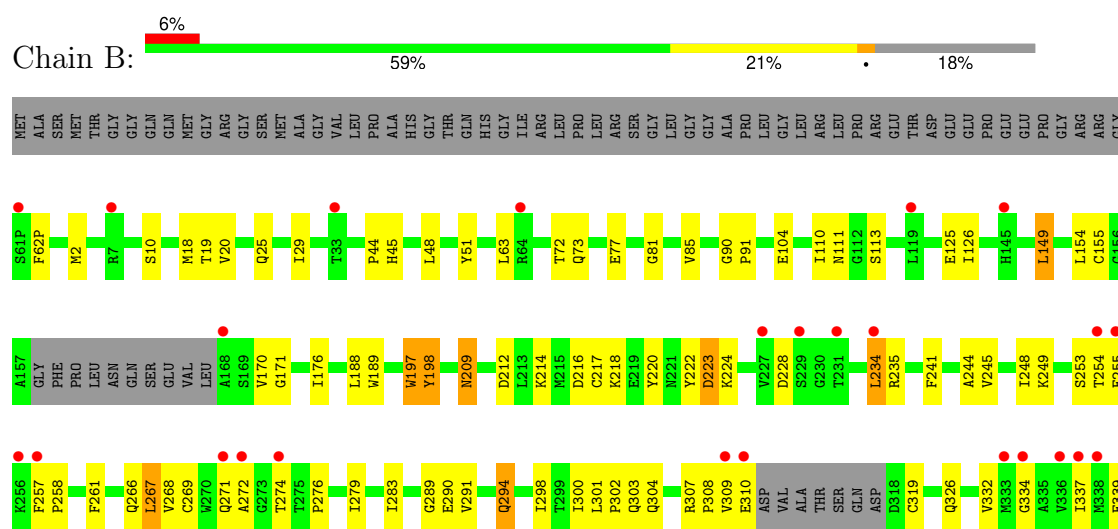
3 Residue-property plots [i](#)

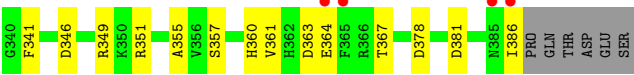
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Beta-secretase 1

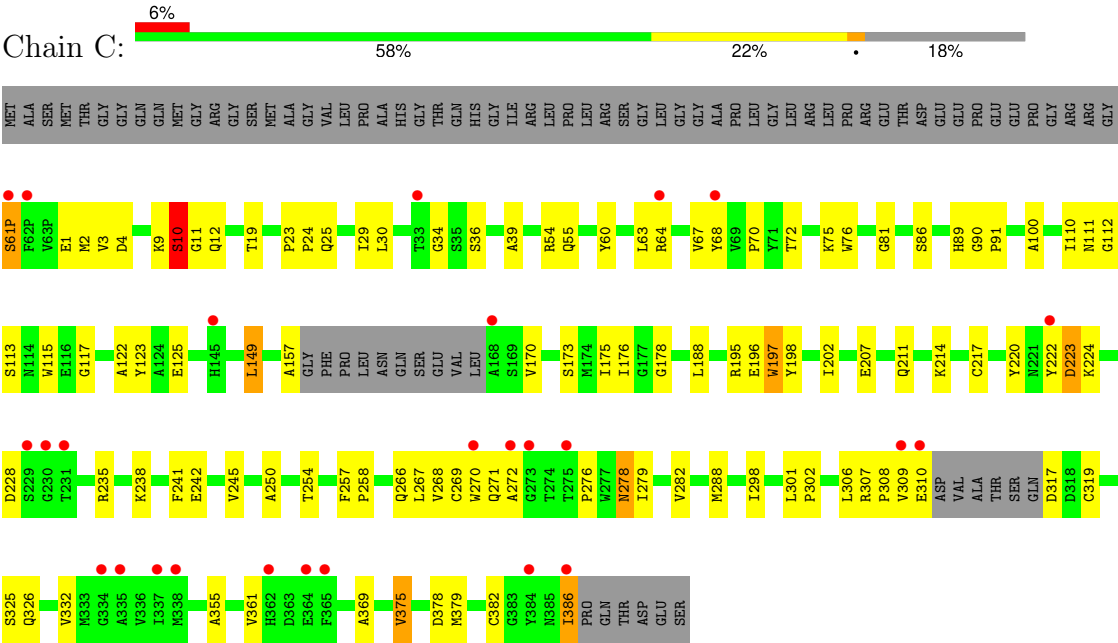


• Molecule 1: Beta-secretase 1





● Molecule 1: Beta-secretase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.96Å 103.35Å 101.01Å 90.00° 104.12° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 31.61 – 2.47	Depositor EDS
% Data completeness (in resolution range)	83.7 (20.00-2.50) 82.3 (31.61-2.47)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.48Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.207 , 0.242 0.198 , 0.232	Depositor DCC
R_{free} test set	5104 reflections (10.08%)	wwPDB-VP
Wilson B-factor (Å ²)	46.0	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9104	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MR0, PO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.39	0/2998	0.68	1/4072 (0.0%)
1	B	0.37	0/3003	0.68	1/4079 (0.0%)
1	C	0.37	0/3011	0.69	0/4090
All	All	0.38	0/9012	0.68	2/12241 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	193	ILE	N-CA-C	-5.12	97.18	111.00
1	B	198	TYR	N-CA-C	-5.02	97.43	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	2924	0	2844	86	0
1	B	2929	0	2854	75	0
1	C	2937	0	2858	87	0
2	A	10	0	0	1	0
2	B	5	0	0	1	0
2	C	5	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	44	0	40	3	0
3	B	44	0	40	1	0
3	C	44	0	40	4	0
4	A	64	0	0	0	0
4	B	52	0	0	0	0
4	C	46	0	0	1	0
All	All	9104	0	8676	238	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 14.

All (238) 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:C:4:ASP:H	1:C:173:SER:HB3	1.27	0.96
1:C:269:CYS:HG	1:C:319:CYS:HG	0.88	0.87
1:C:4:ASP:N	1:C:173:SER:HB3	1.97	0.79
1:C:271:GLN:H	1:C:271:GLN:CD	1.91	0.74
1:A:9:LYS:HG3	1:A:12:GLN:HG3	1.69	0.74
1:A:375:VAL:CG2	1:C:375:VAL:HG22	2.18	0.73
1:A:149:LEU:HD23	1:A:150:PHE:N	2.02	0.73
1:B:276:PRO:O	1:B:279:ILE:HG12	1.89	0.73
1:A:375:VAL:HG22	1:C:375:VAL:HG22	1.73	0.71
1:C:188:LEU:HD23	1:C:355:ALA:HB2	1.73	0.70
1:C:202:ILE:CD1	1:C:379:MET:HG3	2.22	0.70
1:C:235:ARG:HG3	1:C:332:VAL:HB	1.72	0.69
1:B:291:VAL:HB	1:B:294:GLN:HG2	1.74	0.69
1:A:62(P):PHE:HB2	1:A:2:MET:CE	2.24	0.68
1:A:9:LYS:HD2	1:A:12:GLN:OE1	1.93	0.67
1:A:62(P):PHE:HB2	1:A:2:MET:HE2	1.77	0.67
1:C:271:GLN:CD	1:C:271:GLN:N	2.48	0.66
1:C:61(P):SER:N	1:C:175:ILE:CG2	2.60	0.65
1:B:155:CYS:O	1:B:170:VAL:HG13	1.97	0.65
1:B:267:LEU:HD12	1:B:267:LEU:H	1.62	0.64
1:B:241:PHE:O	1:B:245:VAL:HG23	1.98	0.64
1:A:261:PHE:CD1	1:A:268:VAL:HG23	2.33	0.63
1:B:291:VAL:O	1:B:294:GLN:HG3	1.98	0.63
1:A:300:ILE:HG12	1:A:337:ILE:CD1	2.28	0.63
1:C:61(P):SER:N	1:C:175:ILE:HG23	2.14	0.62
1:C:149:LEU:HD23	1:C:149:LEU:O	1.99	0.62
1:B:155:CYS:O	1:B:170:VAL:CG1	2.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:HD11	3:A:801:MR0:H121	1.82	0.60
1:C:68:TYR:OH	1:C:75:LYS:HD3	2.01	0.60
1:B:235:ARG:HG3	1:B:332:VAL:HB	1.84	0.60
1:B:244:ALA:O	1:B:248:ILE:HG13	2.01	0.60
1:A:188:LEU:HD23	1:A:355:ALA:HB2	1.84	0.59
1:A:235:ARG:HG3	1:A:332:VAL:HB	1.84	0.59
1:B:188:LEU:HD23	1:B:355:ALA:HB2	1.84	0.59
1:A:269:CYS:HG	1:A:319:CYS:CB	2.15	0.59
1:A:276:PRO:O	1:A:279:ILE:HG12	2.03	0.59
1:C:235:ARG:CG	1:C:332:VAL:HB	2.33	0.58
1:C:267:LEU:HD22	1:C:309:VAL:HG23	1.84	0.58
1:B:258:PRO:HG3	1:B:266:GLN:NE2	2.18	0.58
1:C:301:LEU:HB3	1:C:302:PRO:HD2	1.86	0.58
1:A:222:TYR:O	1:A:223:ASP:CB	2.51	0.57
1:C:288:MET:CE	1:C:378:ASP:HA	2.33	0.57
1:A:110:ILE:HB	1:A:113:SER:HB3	1.87	0.57
1:C:278:ASN:HD22	1:C:278:ASN:H	1.53	0.57
1:A:235:ARG:NH2	2:A:703:PO4:O2	2.37	0.57
1:B:245:VAL:HG12	1:B:249:LYS:HE3	1.87	0.57
1:B:216:ASP:OD1	1:B:218:LYS:HB2	2.03	0.57
1:A:62(P):PHE:CB	1:A:2:MET:HE2	2.35	0.56
1:C:157:ALA:HB2	1:C:170:VAL:HG12	1.87	0.56
1:A:375:VAL:HG23	1:C:375:VAL:HG22	1.87	0.56
1:B:301:LEU:HB3	1:B:302:PRO:HD2	1.87	0.56
1:A:155:CYS:O	1:A:170:VAL:CG1	2.53	0.56
1:C:19:THR:OG1	1:C:86:SER:HB2	2.05	0.56
1:A:375:VAL:HG22	1:C:375:VAL:HG13	1.88	0.55
1:A:378:ASP:HB3	1:A:381:ASP:OD2	2.06	0.55
1:C:238:LYS:HG3	1:C:326:GLN:OE1	2.05	0.55
1:A:3:VAL:O	1:A:173:SER:CB	2.54	0.55
1:A:249:LYS:HE2	1:A:262:TRP:CD1	2.42	0.55
1:A:61(P):SER:O	1:A:62(P):PHE:HB2	2.06	0.55
1:A:3:VAL:O	1:A:173:SER:HB3	2.07	0.55
1:C:19:THR:HA	1:C:25:GLN:O	2.07	0.54
1:B:212:ASP:O	1:B:214:LYS:HG3	2.07	0.54
1:A:62(P):PHE:CB	1:A:2:MET:CE	2.85	0.54
1:B:170:VAL:HG12	1:B:171:GLY:N	2.22	0.54
1:C:235:ARG:HB2	1:C:332:VAL:HB	1.90	0.54
1:B:253:SER:O	1:B:255:GLU:N	2.40	0.53
1:A:300:ILE:HG12	1:A:337:ILE:HD12	1.88	0.53
1:C:202:ILE:HD13	1:C:379:MET:HG3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:MET:HG2	1:B:90:GLY:HA2	1.89	0.53
1:C:238:LYS:HE2	1:C:242:GLU:OE2	2.08	0.53
1:B:267:LEU:HD12	1:B:267:LEU:N	2.23	0.52
1:A:367:THR:H	1:C:211:GLN:HE22	1.58	0.52
1:B:269:CYS:HG	1:B:319:CYS:CB	2.20	0.52
1:C:63:LEU:HG	1:C:81:GLY:HA2	1.92	0.52
1:C:110:ILE:HB	1:C:113:SER:HB3	1.91	0.52
1:C:241:PHE:CG	1:C:326:GLN:HB3	2.43	0.52
1:C:125:GLU:CD	1:C:195:ARG:HH11	2.13	0.52
1:A:287:LEU:O	1:A:295:SER:HB2	2.10	0.52
1:C:222:TYR:O	1:C:223:ASP:CB	2.57	0.52
1:A:258:PRO:HG3	1:A:266:GLN:NE2	2.25	0.52
1:C:198:TYR:CE2	1:C:224:LYS:HE3	2.45	0.52
1:B:300:ILE:HA	1:B:304:GLN:OE1	2.10	0.52
1:B:303:GLN:HB2	1:B:361:VAL:CG1	2.39	0.51
1:B:125:GLU:HG2	1:B:197:TRP:HB3	1.92	0.51
1:A:3:VAL:O	1:A:173:SER:OG	2.28	0.51
1:C:270:TRP:O	1:C:317:ASP:HB3	2.11	0.51
1:B:378:ASP:HB3	1:B:381:ASP:OD2	2.11	0.51
1:B:228:ASP:O	1:B:334:GLY:HA2	2.11	0.50
1:A:42:ALA:CB	1:A:101:ALA:HB1	2.41	0.50
1:A:302:PRO:HA	1:A:305:TYR:CE2	2.47	0.50
1:B:301:LEU:HD11	1:B:367:THR:HA	1.92	0.50
1:C:34:GLY:O	3:C:803:MR0:H25	2.12	0.50
1:A:217:CYS:HG	1:A:382:CYS:HG	0.54	0.50
1:A:155:CYS:HG	1:A:359:CYS:HG	0.62	0.49
1:B:126:ILE:HG23	1:B:197:TRP:HB2	1.93	0.49
1:C:288:MET:HE3	1:C:378:ASP:HA	1.94	0.49
1:A:205:ARG:HB3	1:A:286:TYR:HB2	1.94	0.49
1:B:91:PRO:HD3	1:B:176:ILE:HB	1.95	0.49
1:B:72:THR:HB	3:B:802:MR0:H3	1.94	0.49
1:C:29:ILE:HD12	1:C:117:GLY:HA3	1.94	0.49
1:B:364:GLU:N	1:B:364:GLU:CD	2.66	0.49
1:A:291:VAL:CG1	1:B:189:TRP:CZ3	2.95	0.49
1:B:290:GLU:N	1:B:294:GLN:OE1	2.37	0.49
1:B:271:GLN:CD	1:B:271:GLN:H	2.16	0.49
1:B:289:GLY:HA3	1:B:294:GLN:O	2.12	0.49
1:A:194:ARG:HD3	1:A:200:GLU:OE2	2.13	0.49
1:A:278:ASN:OD1	1:C:254:THR:HG23	2.13	0.49
1:B:77:GLU:HG2	1:B:104:GLU:HB2	1.95	0.49
1:C:36:SER:OG	1:C:122:ALA:HB3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:ASP:HB3	1:A:366:ARG:O	2.13	0.48
1:A:170:VAL:HG12	1:A:171:GLY:N	2.28	0.48
1:B:245:VAL:CG1	1:B:249:LYS:HE3	2.43	0.48
1:A:63:LEU:HG	1:A:81:GLY:HA2	1.95	0.48
1:A:235:ARG:HB2	1:A:332:VAL:HB	1.95	0.48
1:B:198:TYR:CE2	1:B:224:LYS:HE3	2.49	0.48
1:C:2:MET:HG2	1:C:90:GLY:HA2	1.95	0.48
1:B:267:LEU:HD23	1:B:309:VAL:HG21	1.95	0.48
1:B:217:CYS:HA	1:B:220:TYR:CD1	2.48	0.48
1:B:302:PRO:HG2	1:B:303:GLN:OE1	2.13	0.48
1:C:72:THR:HB	3:C:803:MR0:H3	1.95	0.48
1:C:302:PRO:O	1:C:306:LEU:HB2	2.14	0.48
1:A:359:CYS:O	1:A:359:CYS:SG	2.72	0.47
1:A:365:PHE:CG	1:C:250:ALA:HB1	2.49	0.47
1:C:307:ARG:HA	1:C:308:PRO:HD3	1.72	0.47
1:A:63(P):VAL:HG13	1:A:1:GLU:N	2.28	0.47
1:B:222:TYR:O	1:B:223:ASP:HB3	2.14	0.47
1:B:20:VAL:HG12	1:B:85:VAL:HG22	1.96	0.47
1:C:235:ARG:NH2	2:C:701:PO4:O2	2.30	0.47
1:A:110:ILE:HD11	3:A:801:MR0:H101	1.97	0.47
1:A:300:ILE:HG23	1:A:304:GLN:HB2	1.97	0.47
1:B:149:LEU:C	1:B:149:LEU:HD23	2.34	0.47
1:C:125:GLU:O	1:C:125:GLU:HG3	2.13	0.47
1:C:298:ILE:O	1:C:298:ILE:HG13	2.15	0.47
1:A:9:LYS:HG3	1:A:12:GLN:CG	2.42	0.47
1:B:44:PRO:HD3	1:B:51:TYR:CZ	2.50	0.47
1:B:298:ILE:O	1:B:298:ILE:HG13	2.15	0.47
1:B:19:THR:HA	1:B:25:GLN:O	2.15	0.47
1:C:10:SER:HB3	1:C:11:GLY:H	1.49	0.47
1:C:267:LEU:O	1:C:267:LEU:HD12	2.14	0.47
1:C:197:TRP:CG	1:C:198:TYR:N	2.84	0.46
1:A:149:LEU:HD23	1:A:149:LEU:C	2.35	0.46
1:A:170:VAL:HG12	1:A:171:GLY:H	1.79	0.46
1:A:179:ILE:HD11	1:A:344:VAL:HG11	1.97	0.46
1:B:357:SER:O	1:B:360:HIS:HB3	2.16	0.46
1:C:12:GLN:OE1	1:C:113:SER:HA	2.15	0.46
1:B:222:TYR:O	1:B:223:ASP:CB	2.63	0.46
1:A:207:GLU:HG2	1:A:212:ASP:HA	1.97	0.46
1:B:257:PHE:HD2	1:B:268:VAL:HG21	1.80	0.46
1:B:267:LEU:HD23	1:B:309:VAL:CG2	2.46	0.46
1:B:110:ILE:HB	1:B:113:SER:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:ILE:HB	1:B:300:ILE:HG12	1.98	0.46
1:C:238:LYS:O	1:C:242:GLU:HG3	2.15	0.46
1:B:234:LEU:HG	1:B:337:ILE:HD11	1.98	0.46
1:B:386:ILE:H	1:B:386:ILE:HG13	1.44	0.45
1:C:282:VAL:HG12	1:C:301:LEU:HD23	1.98	0.45
1:C:9:LYS:HE2	1:C:12:GLN:OE1	2.16	0.45
1:C:217:CYS:HA	1:C:220:TYR:CD1	2.51	0.45
1:A:217:CYS:HA	1:A:220:TYR:CD1	2.51	0.45
1:A:344:VAL:O	1:A:352:ILE:HA	2.17	0.45
1:C:149:LEU:HD21	1:C:178:GLY:HA2	1.97	0.45
1:B:209:ASN:HD22	1:B:209:ASN:HA	1.62	0.45
1:C:241:PHE:O	1:C:245:VAL:HG23	2.17	0.45
1:A:29:ILE:HD12	1:A:117:GLY:HA3	1.98	0.45
1:B:303:GLN:HB2	1:B:361:VAL:HG11	1.98	0.45
1:C:110:ILE:HD11	3:C:803:MR0:H101	1.98	0.45
1:B:271:GLN:O	1:B:272:ALA:C	2.54	0.45
1:B:346:ASP:HB3	1:B:351:ARG:HG3	1.98	0.45
1:A:189:TRP:CZ3	1:B:291:VAL:CG1	3.00	0.44
1:B:261:PHE:CD1	1:B:268:VAL:HG23	2.52	0.44
1:C:235:ARG:CB	1:C:332:VAL:HB	2.48	0.44
1:A:267:LEU:HD23	1:A:309:VAL:HG21	2.00	0.44
1:A:28:ASN:HB2	1:A:115:TRP:HA	1.98	0.44
1:B:294:GLN:HE21	1:B:294:GLN:HB2	1.63	0.44
1:C:214:LYS:HE2	1:C:214:LYS:HB3	1.81	0.44
1:C:369:ALA:HB1	4:C:952:HOH:O	2.18	0.44
1:A:155:CYS:CB	1:A:359:CYS:SG	3.05	0.44
1:A:174:MET:CE	1:A:176:ILE:HD11	2.48	0.44
1:C:30:LEU:HD23	1:C:30:LEU:C	2.38	0.44
1:C:34:GLY:HA3	1:C:228:ASP:OD1	2.17	0.44
1:C:207:GLU:HA	1:C:211:GLN:O	2.18	0.44
1:A:113:SER:O	1:A:114:ASN:HB3	2.17	0.44
1:A:30:LEU:CD2	1:A:32:ASP:HB2	2.48	0.43
1:A:197:TRP:CG	1:A:198:TYR:N	2.86	0.43
1:A:238:LYS:O	1:A:242:GLU:HG3	2.18	0.43
1:B:271:GLN:HB2	1:B:274:THR:HG21	2.01	0.43
1:A:76:TRP:HA	1:A:105:SER:HA	1.99	0.43
1:B:170:VAL:CG1	1:B:171:GLY:N	2.81	0.43
1:C:54:ARG:HD2	1:C:60:TYR:CZ	2.53	0.43
1:C:257:PHE:HD2	1:C:268:VAL:HG21	1.82	0.43
1:A:268:VAL:HG12	1:A:269:CYS:N	2.34	0.43
1:C:91:PRO:HD3	1:C:176:ILE:HB	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:GLY:O	3:A:801:MR0:H122	2.19	0.43
1:A:280:PHE:HB3	1:A:302:PRO:HB3	2.00	0.43
1:C:39:ALA:HB2	1:C:100:ALA:HB3	2.01	0.43
1:B:63:LEU:HG	1:B:81:GLY:HA2	2.00	0.43
1:B:45:HIS:HB3	1:B:48:LEU:HG	2.01	0.43
1:C:67:VAL:HG13	1:C:76:TRP:HZ3	1.84	0.43
1:A:48:LEU:HD21	1:A:109:PHE:CE2	2.53	0.42
1:B:268:VAL:HG12	1:B:269:CYS:N	2.33	0.42
1:A:23:PRO:HA	1:A:24:PRO:HD3	1.82	0.42
1:A:254:THR:HG22	1:A:255:GLU:HG3	2.01	0.42
1:C:9:LYS:O	1:C:10:SER:C	2.57	0.42
1:C:217:CYS:HG	1:C:382:CYS:CB	2.26	0.42
1:C:12:GLN:C	3:C:803:MR0:H111	2.40	0.42
1:B:154:LEU:O	1:B:339:GLU:HA	2.19	0.42
1:C:23:PRO:HA	1:C:24:PRO:HD3	1.88	0.42
1:C:267:LEU:HD12	1:C:267:LEU:C	2.40	0.42
1:B:18:MET:SD	1:B:29:ILE:HG13	2.60	0.42
1:C:30:LEU:HD12	1:C:115:TRP:CE2	2.54	0.42
1:A:271:GLN:O	1:A:274:THR:HG23	2.20	0.42
1:C:3:VAL:O	1:C:4:ASP:HB3	2.20	0.42
1:A:2:MET:HG2	1:A:90:GLY:HA2	2.02	0.41
1:B:241:PHE:CG	1:B:326:GLN:HB3	2.55	0.41
1:C:149:LEU:HD23	1:C:149:LEU:C	2.39	0.41
1:C:386:ILE:O	1:C:386:ILE:HG22	2.20	0.41
1:A:365:PHE:CD1	1:A:366:ARG:HG3	2.54	0.41
1:C:276:PRO:O	1:C:279:ILE:HG12	2.20	0.41
1:A:174:MET:HE3	1:A:176:ILE:HD11	2.01	0.41
1:C:2:MET:HG2	1:C:89:HIS:O	2.19	0.41
1:A:62(P):PHE:CB	1:A:2:MET:HE3	2.49	0.41
1:B:235:ARG:NH2	2:B:704:PO4:O1	2.53	0.41
1:A:48:LEU:HD21	1:A:109:PHE:CD2	2.55	0.41
1:A:349:ARG:HH21	1:B:349:ARG:HH21	1.67	0.41
1:B:307:ARG:HA	1:B:308:PRO:HD3	1.74	0.41
1:A:209:ASN:ND2	1:A:281:PRO:HB3	2.36	0.41
1:B:303:GLN:NE2	1:B:363:ASP:HB3	2.35	0.41
1:B:309:VAL:O	1:B:310:GLU:C	2.59	0.41
1:B:341:PHE:HB3	1:B:355:ALA:O	2.21	0.41
1:C:123:TYR:CZ	1:C:196:GLU:HG2	2.54	0.41
1:A:235:ARG:CG	1:A:332:VAL:HB	2.50	0.41
1:A:241:PHE:O	1:A:245:VAL:HG23	2.21	0.41
1:C:258:PRO:HB2	1:C:266:GLN:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:THR:O	1:A:336:VAL:HG13	2.20	0.40
1:A:304:GLN:O	1:A:336:VAL:HB	2.21	0.40
1:A:268:VAL:O	1:A:319:CYS:HA	2.22	0.40
1:C:241:PHE:CD2	1:C:326:GLN:HB3	2.56	0.40
1:C:9:LYS:HE2	1:C:112:GLY:O	2.21	0.40
1:C:310:GLU:O	1:C:310:GLU:HG3	2.21	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	365/455 (80%)	340 (93%)	21 (6%)	4 (1%)	14	26
1	B	366/455 (80%)	336 (92%)	26 (7%)	4 (1%)	14	26
1	C	367/455 (81%)	345 (94%)	18 (5%)	4 (1%)	14	26
All	All	1098/1365 (80%)	1021 (93%)	65 (6%)	12 (1%)	14	26

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62(P)	PHE
1	A	223	ASP
1	B	10	SER
1	B	223	ASP
1	C	10	SER
1	C	223	ASP
1	B	254	THR
1	C	272	ALA
1	B	73	GLN
1	A	131	ASP
1	A	70	PRO

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Mol	Chain	Res	Type
1	C	70	PRO

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	317/381 (83%)	310 (98%)	7 (2%)	52	77
1	B	317/381 (83%)	309 (98%)	8 (2%)	47	73
1	C	318/381 (84%)	305 (96%)	13 (4%)	30	55
All	All	952/1143 (83%)	924 (97%)	28 (3%)	42	69

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62(P)	PHE
1	A	111	ASN
1	A	173	SER
1	A	197	TRP
1	A	234	LEU
1	A	254	THR
1	A	267	LEU
1	B	62(P)	PHE
1	B	111	ASN
1	B	149	LEU
1	B	197	TRP
1	B	209	ASN
1	B	234	LEU
1	B	267	LEU
1	B	294	GLN
1	C	61(P)	SER
1	C	1	GLU
1	C	10	SER
1	C	55	GLN
1	C	64	ARG
1	C	111	ASN

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Mol	Chain	Res	Type
1	C	149	LEU
1	C	197	TRP
1	C	278	ASN
1	C	325	SER
1	C	361	VAL
1	C	375	VAL
1	C	386	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	98	ASN
1	A	111	ASN
1	A	266	GLN
1	A	271	GLN
1	B	73	GLN
1	B	92	ASN
1	B	98	ASN
1	B	111	ASN
1	B	209	ASN
1	B	266	GLN
1	C	98	ASN
1	C	111	ASN
1	C	211	GLN
1	C	278	ASN
1	C	294	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

7 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	PO4	A	702	-	4,4,4	2.06	3 (75%)	6,6,6	0.43	0
3	MR0	B	802	-	43,46,46	2.49	13 (30%)	44,62,62	2.24	12 (27%)
2	PO4	C	701	-	4,4,4	1.93	3 (75%)	6,6,6	0.46	0
2	PO4	B	704	-	4,4,4	1.42	0	6,6,6	0.50	0
2	PO4	A	703	-	4,4,4	1.38	0	6,6,6	0.45	0
3	MR0	C	803	-	43,46,46	2.25	12 (27%)	44,62,62	1.96	9 (20%)
3	MR0	A	801	-	43,46,46	2.30	14 (32%)	44,62,62	2.00	12 (27%)

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	MR0	B	802	-	-	17/41/42/42	0/3/3/3
3	MR0	C	803	-	-	19/41/42/42	0/3/3/3
3	MR0	A	801	-	-	17/41/42/42	0/3/3/3

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	802	MR0	C25-C20	7.33	1.49	1.36
3	B	802	MR0	C29-C28	6.71	1.48	1.36
3	A	801	MR0	C25-C20	6.35	1.47	1.36
3	B	802	MR0	C6-N	6.21	1.45	1.34
3	C	803	MR0	C25-C20	6.13	1.47	1.36
3	C	803	MR0	C6-N	6.08	1.45	1.34
3	A	801	MR0	C6-N	5.98	1.45	1.34
3	A	801	MR0	C29-C28	5.80	1.46	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	803	MR0	C29-C28	5.75	1.46	1.36
3	B	802	MR0	C8-N1	3.95	1.42	1.34
3	B	802	MR0	C17-C28	3.92	1.58	1.50
3	A	801	MR0	C22-C21	3.46	1.45	1.37
3	A	801	MR0	C15-N1	3.36	1.51	1.46
3	A	801	MR0	C5-C	3.28	1.46	1.38
3	C	803	MR0	C22-C21	3.17	1.44	1.37
3	C	803	MR0	C5-C	3.16	1.46	1.38
3	A	801	MR0	C17-C28	3.14	1.56	1.50
3	C	803	MR0	C17-C28	3.13	1.56	1.50
3	B	802	MR0	C5-C	3.09	1.46	1.38
3	C	803	MR0	C8-N1	3.06	1.40	1.34
3	B	802	MR0	C22-C21	3.03	1.44	1.37
3	B	802	MR0	C17-C15	2.97	1.56	1.53
3	B	802	MR0	C1-C2	2.97	1.45	1.38
3	B	802	MR0	C15-N1	2.96	1.51	1.46
3	C	803	MR0	C15-N1	2.92	1.51	1.46
3	A	801	MR0	C1-C2	2.90	1.45	1.38
3	A	801	MR0	C8-N1	2.87	1.40	1.34
3	A	801	MR0	C10-N	2.67	1.54	1.47
3	C	803	MR0	C1-C2	2.61	1.45	1.38
3	B	802	MR0	C10-N	2.53	1.53	1.47
3	A	801	MR0	C23-C24	2.49	1.45	1.38
2	A	702	PO4	P-O3	-2.41	1.47	1.54
3	C	803	MR0	C23-C24	2.39	1.44	1.38
3	A	801	MR0	C7-N7A	2.35	1.40	1.33
2	C	701	PO4	P-O2	-2.34	1.47	1.54
2	A	702	PO4	P-O4	-2.31	1.47	1.54
3	C	803	MR0	C10-N	2.30	1.53	1.47
3	C	803	MR0	C17-C15	2.27	1.56	1.53
2	A	702	PO4	P-O2	-2.18	1.48	1.54
3	A	801	MR0	C18-N2	2.17	1.50	1.47
2	C	701	PO4	P-O3	-2.16	1.48	1.54
3	B	802	MR0	F26-C26	-2.15	1.30	1.35
3	A	801	MR0	C17-C15	2.13	1.55	1.53
3	B	802	MR0	C18-N2	2.10	1.50	1.47
2	C	701	PO4	P-O4	-2.04	1.48	1.54

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	802	MR0	O3-C24-C25	8.62	131.30	113.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	MR0	O3-C24-C25	6.84	127.64	113.54
3	C	803	MR0	O3-C24-C25	6.83	127.62	113.54
3	B	802	MR0	O3-C24-C23	-5.65	111.90	124.54
3	C	803	MR0	C17-C15-N1	-4.83	104.28	109.58
3	B	802	MR0	C17-C15-N1	-4.40	104.76	109.58
3	A	801	MR0	C17-C15-N1	-4.39	104.76	109.58
3	A	801	MR0	O3-C24-C23	-4.27	114.97	124.54
3	B	802	MR0	CO-O3-C24	4.21	126.14	117.23
3	C	803	MR0	C16-C15-N1	4.09	117.64	109.97
3	C	803	MR0	O3-C24-C23	-4.03	115.53	124.54
3	A	801	MR0	C16-C15-N1	3.96	117.39	109.97
3	B	802	MR0	C16-C18-N2	-3.17	107.79	111.95
3	A	801	MR0	O-C6-N	-3.05	117.85	122.67
3	B	802	MR0	C16-C15-N1	2.98	115.56	109.97
3	A	801	MR0	C30-C29-C28	-2.96	119.77	124.48
3	C	803	MR0	C30-C29-C28	-2.92	119.83	124.48
3	A	801	MR0	CO-O3-C24	2.77	123.08	117.23
3	A	801	MR0	C1-C2-C7	2.69	121.52	117.86
3	B	802	MR0	C30-C29-C28	-2.61	120.33	124.48
3	B	802	MR0	C25-C24-C23	-2.58	116.80	120.96
3	C	803	MR0	C25-C24-C23	-2.55	116.85	120.96
3	C	803	MR0	CO-O3-C24	2.47	122.44	117.23
3	C	803	MR0	O-C6-N	-2.40	118.88	122.67
3	A	801	MR0	C10-N-C6	2.36	127.19	120.66
3	B	802	MR0	C10-N-C6	2.31	127.05	120.66
3	B	802	MR0	C1-C2-C7	2.29	120.97	117.86
3	C	803	MR0	C16-C18-N2	-2.28	108.96	111.95
3	B	802	MR0	O-C6-N	-2.24	119.13	122.67
3	A	801	MR0	C25-C24-C23	-2.22	117.38	120.96
3	A	801	MR0	C9-N-C6	-2.19	114.60	120.66
3	B	802	MR0	C9-N-C6	-2.17	114.67	120.66
3	A	801	MR0	C13-C10-N	2.04	120.92	112.37

There are no chirality outliers.

All (53) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	801	MR0	C5-C4-C8-O1
3	A	801	MR0	C-C6-N-C10
3	A	801	MR0	C-C6-N-C9
3	A	801	MR0	O-C6-N-C10
3	A	801	MR0	O-C6-N-C9

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Mol	Chain	Res	Type	Atoms
3	A	801	MR0	C13-C10-N-C6
3	B	802	MR0	C3-C4-C8-O1
3	B	802	MR0	C3-C2-C7-N7A
3	B	802	MR0	N2-C19-C20-C25
3	C	803	MR0	C5-C4-C8-N1
3	C	803	MR0	C5-C4-C8-O1
3	C	803	MR0	C-C6-N-C9
3	C	803	MR0	O-C6-N-C9
3	C	803	MR0	C1-C2-C7-N7A
3	C	803	MR0	C3-C2-C7-N7A
3	C	803	MR0	C1-C2-C7-O7B
3	C	803	MR0	C13-C10-N-C6
3	A	801	MR0	C12-C11-C9-N
3	C	803	MR0	C12-C11-C9-N
3	C	803	MR0	O-C6-N-C10
3	B	802	MR0	C12-C11-C9-N
3	B	802	MR0	C13-C10-N-C6
3	B	802	MR0	O-C6-N-C9
3	A	801	MR0	C5-C4-C8-N1
3	B	802	MR0	O-C6-N-C10
3	A	801	MR0	C13-C10-N-C9
3	A	801	MR0	C11-C9-N-C10
3	B	802	MR0	C-C6-N-C9
3	C	803	MR0	C-C6-N-C10
3	C	803	MR0	C13-C10-N-C9
3	B	802	MR0	C3-C2-C7-O7B
3	B	802	MR0	C3-C4-C8-N1
3	A	801	MR0	C5-C-C6-O
3	A	801	MR0	C1-C-C6-O
3	B	802	MR0	C5-C4-C8-O1
3	B	802	MR0	C5-C4-C8-N1
3	C	803	MR0	C3-C4-C8-O1
3	A	801	MR0	C25-C24-O3-CO
3	B	802	MR0	C25-C24-O3-CO
3	C	803	MR0	C25-C24-O3-CO
3	A	801	MR0	C23-C24-O3-CO
3	B	802	MR0	C23-C24-O3-CO
3	C	803	MR0	C23-C24-O3-CO
3	C	803	MR0	C1-C-C6-O
3	A	801	MR0	C11-C9-N-C6
3	C	803	MR0	C11-C9-N-C10
3	B	802	MR0	C-C6-N-C10

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Mol	Chain	Res	Type	Atoms
3	B	802	MR0	C13-C10-N-C9
3	C	803	MR0	C3-C2-C7-O7B
3	B	802	MR0	C1-C-C6-O
3	A	801	MR0	C5-C-C6-N
3	A	801	MR0	C3-C2-C7-O7B
3	C	803	MR0	C11-C9-N-C6

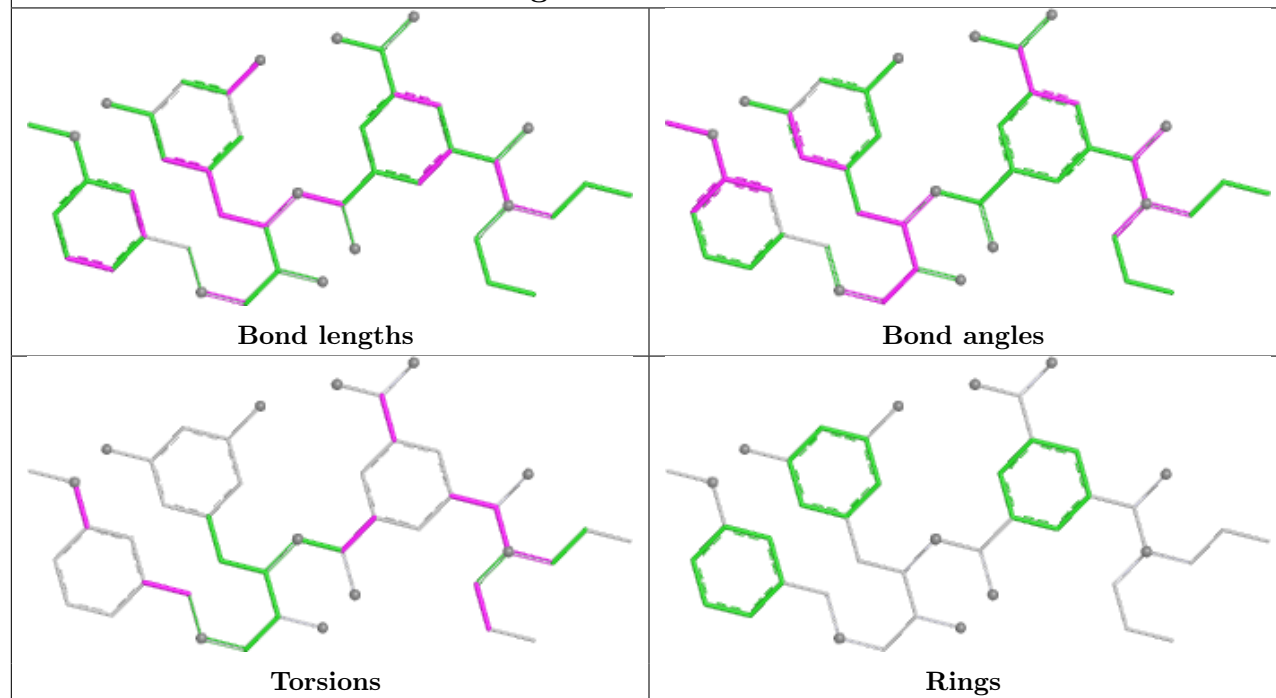
There are no ring outliers.

6 monomers are involved in 11 short contacts:

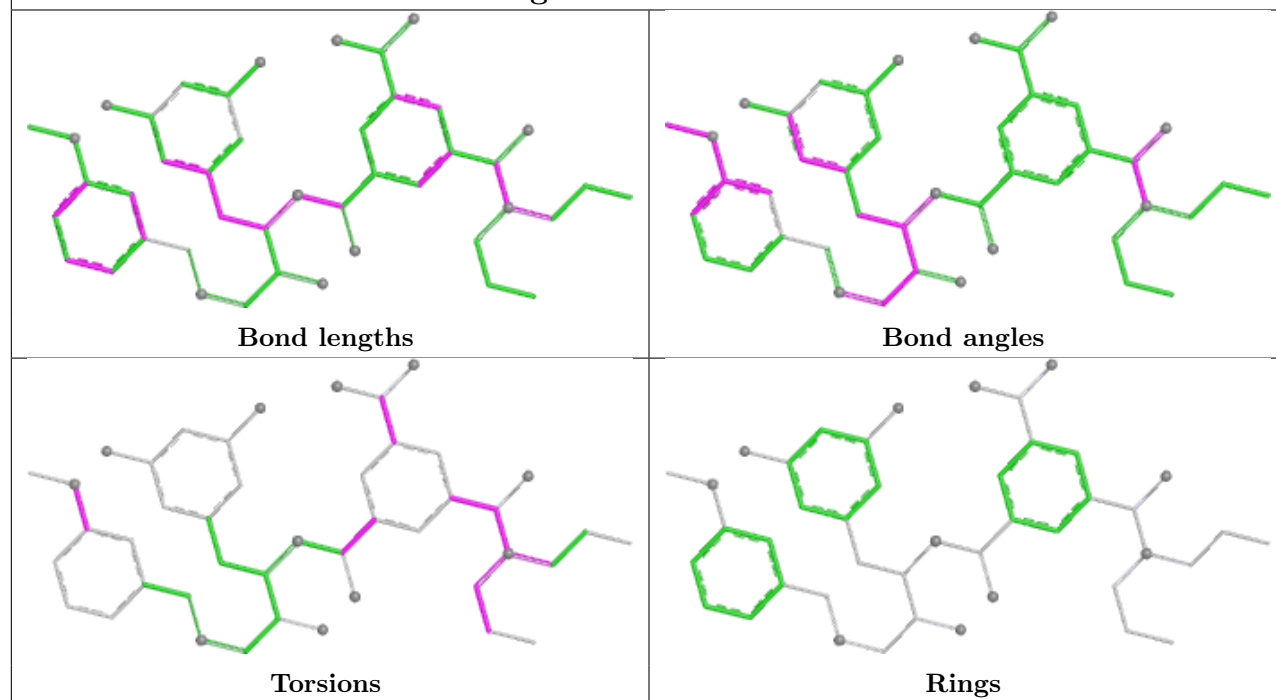
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	802	MR0	1	0
2	C	701	PO4	1	0
2	B	704	PO4	1	0
2	A	703	PO4	1	0
3	C	803	MR0	4	0
3	A	801	MR0	3	0

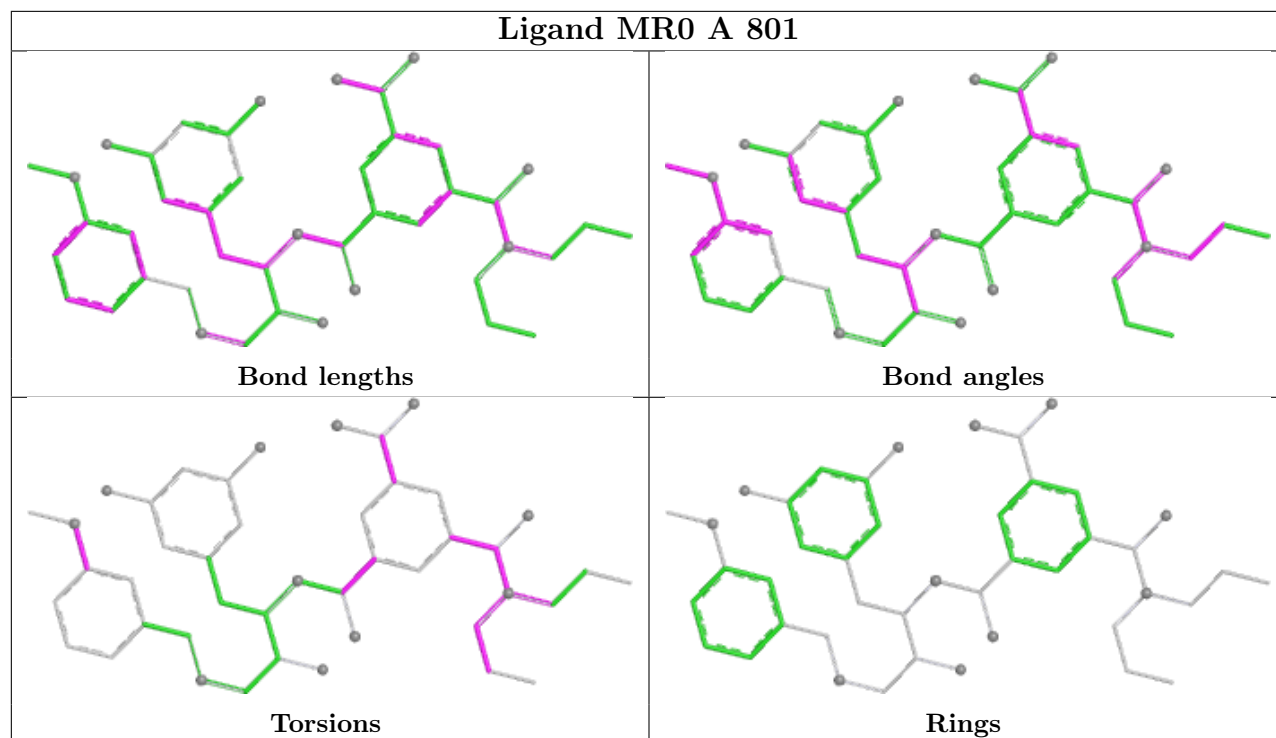
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

Ligand MR0 B 802



Ligand MR0 C 803





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	371/455 (81%)	0.06	15 (4%) 38 41	26, 41, 65, 81	0
1	B	372/455 (81%)	0.32	29 (7%) 13 13	23, 46, 74, 89	0
1	C	373/455 (81%)	0.21	26 (6%) 16 16	26, 44, 70, 86	0
All	All	1116/1365 (81%)	0.20	70 (6%) 20 21	23, 44, 71, 89	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	386	ILE	5.5
1	C	272	ALA	5.3
1	B	168	ALA	5.2
1	B	61(P)	SER	4.6
1	B	310	GLU	4.4
1	C	364	GLU	4.4
1	B	256	LYS	4.3
1	B	386	ILE	3.9
1	A	157	ALA	3.9
1	A	61(P)	SER	3.8
1	B	337	ILE	3.5
1	C	365	PHE	3.5
1	A	362	HIS	3.5
1	B	227	VAL	3.4
1	C	168	ALA	3.4
1	A	227	VAL	3.3
1	B	309	VAL	3.3
1	B	274	THR	3.2
1	B	64	ARG	3.1
1	B	254	THR	3.1
1	C	62(P)	PHE	3.0
1	B	272	ALA	3.0
1	A	364	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	310	GLU	2.9
1	B	229	SER	2.9
1	C	309	VAL	2.9
1	C	229	SER	2.9
1	C	33	THR	2.8
1	B	33	THR	2.8
1	B	385	ASN	2.8
1	C	61(P)	SER	2.8
1	C	64	ARG	2.8
1	B	145	HIS	2.8
1	B	365	PHE	2.7
1	C	275	THR	2.7
1	A	7	ARG	2.6
1	B	255	GLU	2.6
1	B	364	GLU	2.6
1	C	362	HIS	2.6
1	B	7	ARG	2.5
1	A	365	PHE	2.5
1	A	119	LEU	2.4
1	B	234	LEU	2.4
1	C	145	HIS	2.4
1	C	231	THR	2.4
1	A	229	SER	2.4
1	B	333	MET	2.4
1	B	257	PHE	2.4
1	A	121	LEU	2.4
1	C	338	MET	2.3
1	A	169	SER	2.3
1	B	334	GLY	2.3
1	C	270	TRP	2.3
1	C	335	ALA	2.3
1	C	230	GLY	2.3
1	C	384	TYR	2.3
1	B	338	MET	2.2
1	B	271	GLN	2.2
1	C	273	GLY	2.2
1	C	337	ILE	2.2
1	B	119	LEU	2.2
1	A	385	ASN	2.1
1	B	231	THR	2.1
1	C	334	GLY	2.1
1	A	33	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	32	ASP	2.1
1	C	222	TYR	2.1
1	A	62(P)	PHE	2.1
1	B	336	VAL	2.0
1	C	68	TYR	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 monosaccharides in this entry.

6.4 Ligands [i](#)

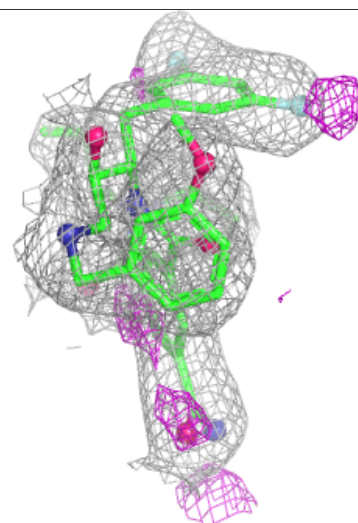
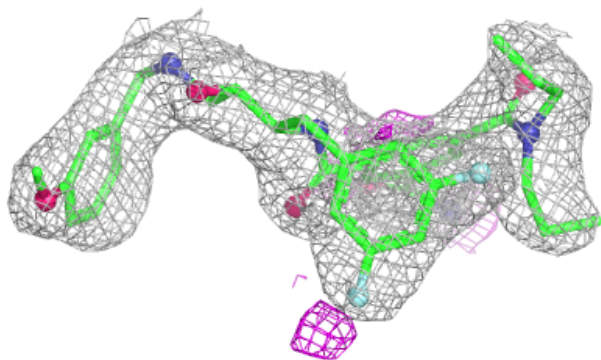
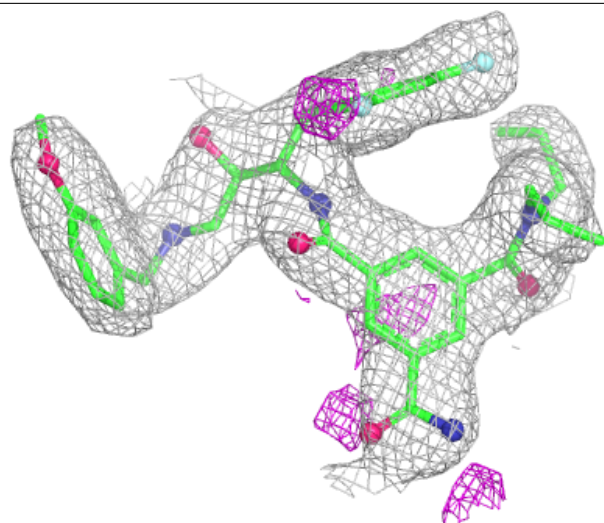
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	B	704	5/5	0.61	0.42	99,99,99,99	0
2	PO4	C	701	5/5	0.73	0.37	99,99,99,99	0
2	PO4	A	703	5/5	0.75	0.38	96,96,97,98	0
3	MR0	B	802	44/44	0.90	0.23	38,43,45,49	0
2	PO4	A	702	5/5	0.91	0.45	99,99,99,99	0
3	MR0	C	803	44/44	0.93	0.22	36,40,45,49	0
3	MR0	A	801	44/44	0.94	0.18	29,35,39,43	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.

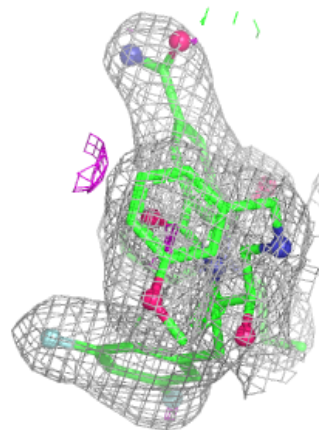
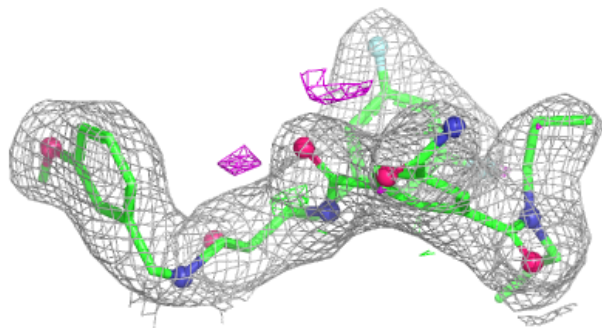
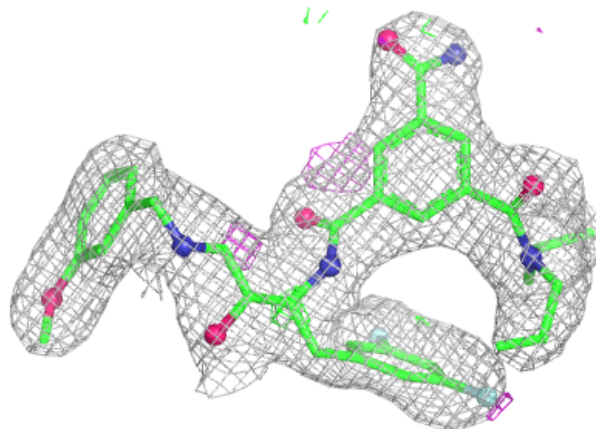
Electron density around MR0 B 802:

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



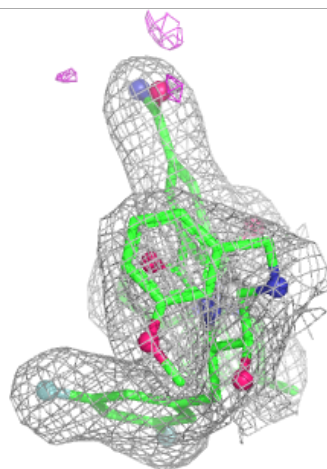
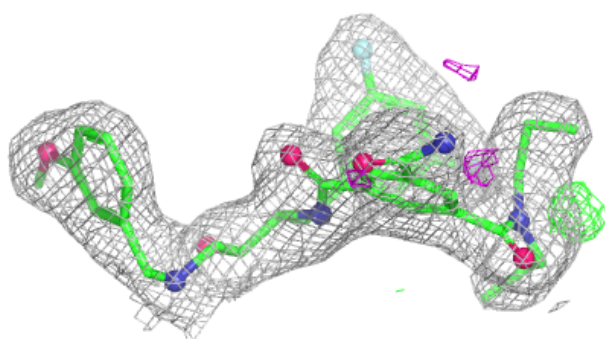
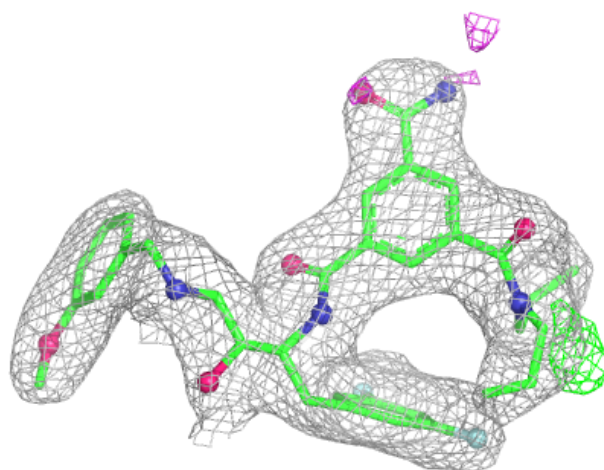
Electron density around MR0 C 803:

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



Electron density around MR0 A 801:

$2mF_o-DF_c$ (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.