



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:10 AM UTC

PDB ID : 2P83 / pdb_00002p83
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 : 4-5-2 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

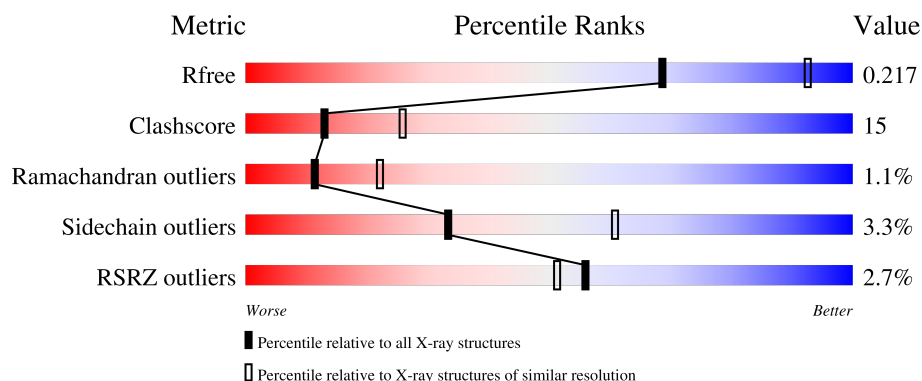
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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	0% 55% 25% • 18%
1	B	455	3% 56% 22% • 18%

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:

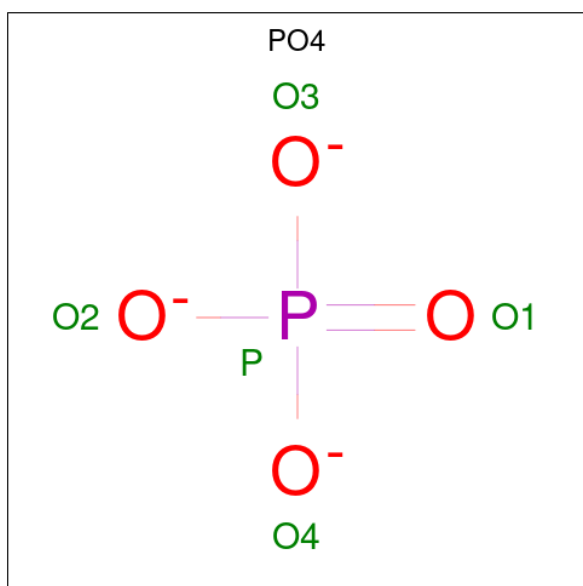
Chain	Residue	Modelled	Actual	Comment	Reference
A	1P	MET	-	expression tag	UNP P56817
A	2P	ALA	-	expression tag	UNP P56817
A	3P	SER	-	expression tag	UNP P56817
A	4P	MET	-	expression tag	UNP P56817
A	5P	THR	-	expression tag	UNP P56817
A	6P	GLY	-	expression tag	UNP P56817
A	7P	GLY	-	expression tag	UNP P56817
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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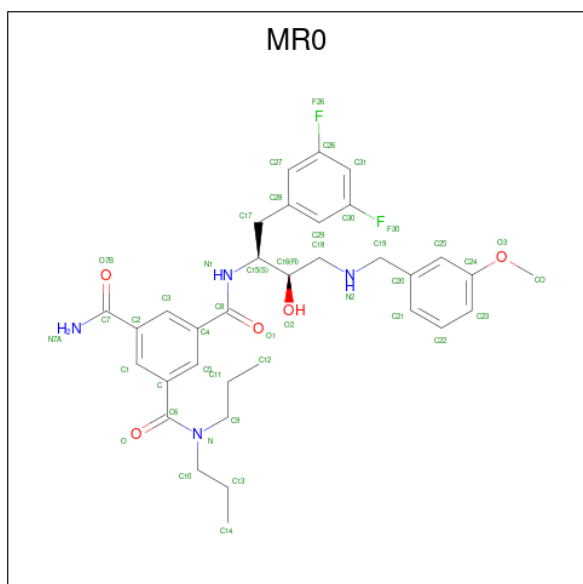
Chain	Residue	Modelled	Actual	Comment	Reference
B	9P	GLN	-	expression tag	UNP P56817
B	10P	MET	-	expression tag	UNP P56817
B	11P	GLY	-	expression tag	UNP P56817
B	12P	ARG	-	expression tag	UNP P56817
B	13P	GLY	-	expression tag	UNP P56817
B	14P	SER	-	expression tag	UNP P56817
B	15P	MET	-	expression tag	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 (CCD ID: PO4) (formula: O₄P).



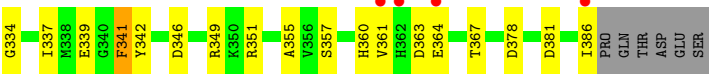
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0

- 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 (CCD ID: MR0) (formula: C₃₃H₄₀F₂N₄O₅).

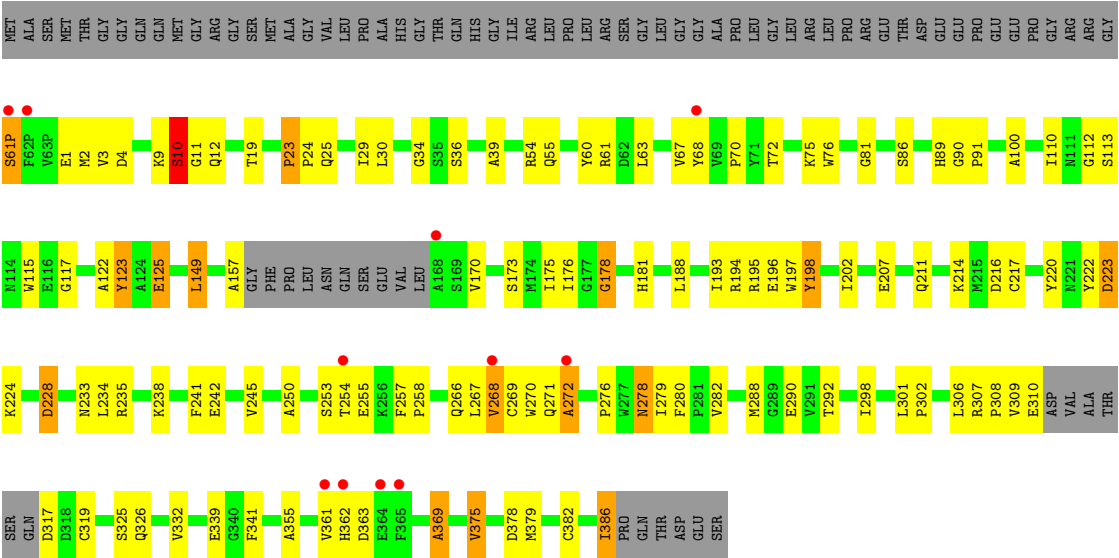


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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	46	Total	O	0	0
			46	46		



● 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 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	83.7 (20.00-2.50) 83.5 (20.00-2.50)	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 (Not available) , 0.217	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)	0.35 , 36.5	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

5.1 Standard geometry

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

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	0.44	1/2998 (0.0%)	1.02	20/4072 (0.5%)
1	B	0.42	0/3003	1.01	11/4079 (0.3%)
1	C	0.43	0/3011	1.05	19/4090 (0.5%)
All	All	0.43	1/9012 (0.0%)	1.03	50/12241 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	3	VAL	CA-CB	-6.07	1.47	1.54

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	198	TYR	N-CA-C	-8.98	98.95	110.53
1	B	198	TYR	N-CA-C	-8.80	97.43	110.46
1	A	3	VAL	N-CA-C	8.60	119.39	110.36
1	C	363	ASP	N-CA-C	-7.79	97.90	110.20
1	C	233	ASN	N-CA-C	7.62	119.79	110.41
1	A	341	PHE	N-CA-C	7.33	121.02	109.81
1	A	233	ASN	N-CA-C	7.20	119.82	110.53
1	A	198	TYR	N-CA-C	-7.10	99.35	110.36
1	A	234	LEU	N-CA-C	-7.09	97.69	108.96
1	B	123	TYR	N-CA-C	7.04	120.06	110.55
1	C	280	PHE	CA-C-N	7.02	126.98	120.03
1	C	280	PHE	C-N-CA	7.02	126.98	120.03
1	B	87	ILE	CA-C-N	6.86	126.37	119.24
1	B	87	ILE	C-N-CA	6.86	126.37	119.24
1	A	193	ILE	N-CA-C	-6.76	97.18	107.73
1	A	307	ARG	CA-C-N	6.53	126.76	119.90
1	A	307	ARG	C-N-CA	6.53	126.76	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	23	PRO	CA-C-N	6.45	127.90	119.84
1	A	23	PRO	C-N-CA	6.45	127.90	119.84
1	C	234	LEU	N-CA-C	-6.39	98.79	108.96
1	C	362	HIS	N-CA-C	6.36	118.42	107.93
1	C	341	PHE	N-CA-C	6.28	118.95	109.41
1	B	72	THR	N-CA-C	-6.23	104.41	111.07
1	A	123	TYR	N-CA-C	6.08	118.76	110.55
1	C	193	ILE	N-CA-C	-6.05	98.29	107.73
1	B	233	ASN	N-CA-C	6.02	118.30	110.53
1	C	339	GLU	N-CA-C	5.99	118.58	111.33
1	C	216	ASP	N-CA-C	-5.99	101.04	109.96
1	C	123	TYR	N-CA-C	5.95	119.33	110.52
1	A	196	GLU	N-CA-C	5.90	118.82	108.56
1	A	72	THR	N-CA-C	-5.87	104.80	111.07
1	A	178	GLY	N-CA-C	5.83	118.73	110.74
1	B	234	LEU	N-CA-C	-5.78	98.33	108.20
1	B	341	PHE	N-CA-C	5.66	117.57	109.14
1	C	290	GLU	N-CA-C	-5.55	105.32	111.71
1	B	342	TYR	N-CA-C	-5.53	99.72	108.73
1	C	23	PRO	CA-C-N	5.42	126.61	119.84
1	C	23	PRO	C-N-CA	5.42	126.61	119.84
1	A	363	ASP	N-CA-C	-5.35	99.40	110.80
1	B	193	ILE	N-CA-C	-5.28	100.00	107.98
1	A	216	ASP	N-CA-C	-5.26	102.13	109.96
1	A	375	VAL	CB-CA-C	-5.25	103.70	110.84
1	B	29	ILE	N-CA-C	5.25	115.66	107.99
1	C	181	HIS	N-CA-C	5.20	118.72	112.38
1	A	369	ALA	N-CA-C	5.12	116.08	108.60
1	A	190	TYR	N-CA-C	5.11	118.13	109.76
1	A	328	SER	N-CA-C	-5.09	105.97	113.61
1	C	228	ASP	N-CA-C	5.06	117.87	107.69
1	C	369	ALA	N-CA-C	5.06	116.50	108.96
1	C	178	GLY	N-CA-C	5.01	118.45	111.19

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	2924	0	2844	89	0
1	B	2929	0	2854	82	0
1	C	2937	0	2858	95	0
2	A	10	0	0	1	0
2	B	5	0	0	1	0
2	C	5	0	0	1	0
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	256	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 (256) 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	0.79
1:C:288:MET:HE1	1:C:379:MET:N	1.99	0.77
1:A:62(P):PHE:HB2	1:A:2:MET:HE2	1.67	0.76
1:A:174:MET:HE3	1:A:176:ILE:HD11	1.68	0.76
1:C:194:ARG:HE	1:C:379:MET:HE3	1.50	0.76
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:62(P):PHE:CB	1:A:2:MET:HE2	2.20	0.72
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:271:GLN:CD	1:C:271:GLN:N	2.48	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:LYS:HD2	1:A:12:GLN:OE1	1.93	0.67
1:A:62(P):PHE:HB2	1:A:2:MET:CE	2.24	0.67
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:174:MET:HE3	1:B:176:ILE:HD11	1.79	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:C:197:TRP:CG	1:C:198:TYR:H	2.16	0.62
1:B:155:CYS:O	1:B:170:VAL:CG1	2.49	0.61
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:197:TRP:CG	1:A:198:TYR:H	2.21	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:288:MET:HE1	1:C:378:ASP:C	2.27	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:C:125:GLU:CD	1:C:195:ARG:HH11	2.13	0.57
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:197:TRP:CD1	1:A:198:TYR:H	2.25	0.54
1:B:212:ASP:O	1:B:214:LYS:HG3	2.07	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:271:GLN:CD	1:B:271:GLN:H	2.16	0.54
1:B:364:GLU:N	1:B:364:GLU:CD	2.66	0.53
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
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:A:62(P):PHE:CB	1:A:2:MET:CE	2.85	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: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:B:149:LEU:C	1:B:149:LEU:HD23	2.34	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:149:LEU:HD23	1:A:149:LEU:C	2.35	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:197:TRP:CG	1:B:198:TYR:H	2.29	0.50
1:B:228:ASP:O	1:B:334:GLY:HA2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:A:291:VAL:CG1	1:B:189:TRP:CZ3	2.95	0.49
1:B:271:GLN:O	1:B:272:ALA:C	2.54	0.49
1:B:290:GLU:N	1:B:294:GLN:OE1	2.37	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
1:C:30:LEU:HD23	1:C:30:LEU:C	2.38	0.48
1:A:363:ASP:HB3	1:A:366:ARG:O	2.13	0.48
1:B:113:SER:HG	1:B:115:TRP:CD1	2.30	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:197:TRP:CD1	1:B:198:TYR:H	2.32	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:C:9:LYS:O	1:C:10:SER:C	2.57	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:12:GLN:C	3:C:803:MR0:H111	2.40	0.47
1:C:298:ILE:O	1:C:298:ILE:HG13	2.15	0.47
1:C:125:GLU:O	1:C:125:GLU:HG3	2.13	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:C:267:LEU:C	1:C:267:LEU:HD12	2.40	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:HD12	1:C:267:LEU:O	2.14	0.47
1:C:197:TRP:CG	1:C:198:TYR:N	2.84	0.46
1:A:170:VAL:HG12	1:A:171:GLY:H	1.79	0.46
1:C:149:LEU:HD23	1:C:149:LEU:C	2.39	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
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:B:309:VAL:O	1:B:310:GLU:C	2.59	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:346:ASP:HB3	1:B:351:ARG:HG3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:253:SER:C	1:B:255:GLU:N	2.75	0.44
1:C:197:TRP:CD1	1:C:198:TYR:H	2.36	0.44
1:B:9:LYS:O	1:B:10:SER:C	2.60	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:271:GLN:O	1:A:272:ALA:C	2.60	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
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:217:CYS:HG	1:C:382:CYS:CB	2.26	0.42
1:C:271:GLN:O	1:C:272:ALA:C	2.61	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:B:18:MET:SD	1:B:29:ILE:HG13	2.60	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:GLY:C	1:B:75:LYS:HG3	2.43	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:288:MET:HE1	1:C:378:ASP:HA	2.01	0.41
1:C:386:ILE:O	1:C:386:ILE:HG22	2.20	0.41
1:A:145:HIS:O	1:A:146:VAL:C	2.62	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:C:2:MET:HG2	1:C:89:HIS:O	2.19	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:C:288:MET:HE1	1:C:378:ASP:CA	2.49	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: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:9:LYS:O	1:A:10:SER:C	2.63	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
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:253:SER:C	1:C:255:GLU:H	2.29	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

5.3.1 Protein backbone

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	365/455 (80%)	340 (93%)	21 (6%)	4 (1%)	11	22
1	B	366/455 (80%)	336 (92%)	26 (7%)	4 (1%)	11	22
1	C	367/455 (81%)	345 (94%)	18 (5%)	4 (1%)	11	22
All	All	1098/1365 (80%)	1021 (93%)	65 (6%)	12 (1%)	11	22

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
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%)	45	73
1	B	317/381 (83%)	307 (97%)	10 (3%)	34	62
1	C	318/381 (84%)	304 (96%)	14 (4%)	25	50
All	All	952/1143 (83%)	921 (97%)	31 (3%)	33	61

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62(P)	PHE
1	A	62	ASP
1	A	173	SER
1	A	234	LEU
1	A	254	THR
1	A	361	VAL
1	A	375	VAL
1	B	62(P)	PHE
1	B	62	ASP
1	B	125	GLU
1	B	149	LEU
1	B	197	TRP
1	B	209	ASN
1	B	213	LEU
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	61	ARG
1	C	125	GLU
1	C	149	LEU
1	C	268	VAL
1	C	278	ASN
1	C	292	THR
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 (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	HIS
1	A	53	GLN
1	A	55	GLN
1	A	98	ASN
1	A	111	ASN
1	A	209	ASN

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Mol	Chain	Res	Type
1	A	211	GLN
1	A	266	GLN
1	A	271	GLN
1	A	385	ASN
1	B	73	GLN
1	B	89	HIS
1	B	92	ASN
1	B	98	ASN
1	B	111	ASN
1	B	209	ASN
1	B	266	GLN
1	B	278	ASN
1	C	53	GLN
1	C	111	ASN
1	C	211	GLN
1	C	278	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	C	701	-	4,4,4	1.93	3 (75%)	6,6,6	0.46	0
2	PO4	A	702	-	4,4,4	2.06	3 (75%)	6,6,6	0.43	0
3	MR0	A	801	-	46,46,46	2.77	26 (56%)	61,62,62	1.77	14 (22%)
3	MR0	C	803	-	46,46,46	2.70	27 (58%)	61,62,62	1.76	14 (22%)
2	PO4	B	704	-	4,4,4	1.42	0	6,6,6	0.50	0
3	MR0	B	802	-	46,46,46	2.87	26 (56%)	61,62,62	1.88	16 (26%)
2	PO4	A	703	-	4,4,4	1.38	0	6,6,6	0.45	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
3	MR0	B	802	-	-	9/42/42/42	0/3/3/3
3	MR0	A	801	-	-	9/42/42/42	0/3/3/3
3	MR0	C	803	-	-	15/42/42/42	0/3/3/3

All (85) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	802	MR0	C25-C20	6.02	1.49	1.39
3	B	802	MR0	C29-C28	5.34	1.48	1.39
3	B	802	MR0	C29-C30	5.28	1.46	1.37
3	B	802	MR0	C6-N	5.11	1.45	1.34
3	C	803	MR0	C6-N	5.00	1.45	1.34
3	A	801	MR0	C25-C20	4.96	1.47	1.39
3	A	801	MR0	C6-N	4.92	1.45	1.34
3	A	801	MR0	C5-C	4.79	1.46	1.39
3	A	801	MR0	C29-C30	4.78	1.45	1.37
3	A	801	MR0	C31-C30	4.77	1.45	1.37
3	B	802	MR0	C31-C30	4.74	1.45	1.37
3	C	803	MR0	C25-C20	4.71	1.47	1.39
3	C	803	MR0	C31-C30	4.69	1.45	1.37
3	C	803	MR0	C5-C	4.60	1.46	1.39
3	C	803	MR0	C29-C30	4.55	1.45	1.37
3	B	802	MR0	C5-C	4.47	1.46	1.39
3	A	801	MR0	C29-C28	4.35	1.46	1.39
3	C	803	MR0	C29-C28	4.30	1.46	1.39
3	B	802	MR0	C1-C2	4.27	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	803	MR0	C22-C23	4.19	1.46	1.38
3	A	801	MR0	C1-C2	4.15	1.45	1.39
3	A	801	MR0	C7-N7A	3.92	1.40	1.33
3	A	801	MR0	C22-C21	3.79	1.45	1.38
3	A	801	MR0	C22-C23	3.76	1.45	1.38
3	B	802	MR0	C3-C2	3.73	1.45	1.39
3	C	803	MR0	C1-C2	3.67	1.45	1.39
3	B	802	MR0	C22-C23	3.59	1.45	1.38
3	C	803	MR0	C5-C4	3.59	1.44	1.39
3	A	801	MR0	C27-C26	3.58	1.43	1.37
3	B	802	MR0	C8-N1	3.55	1.42	1.34
3	C	803	MR0	C31-C26	3.47	1.43	1.37
3	A	801	MR0	C23-C24	3.45	1.45	1.38
3	A	801	MR0	C31-C26	3.44	1.43	1.37
3	C	803	MR0	C22-C21	3.41	1.44	1.38
3	B	802	MR0	C31-C26	3.39	1.43	1.37
3	A	801	MR0	C15-N1	3.36	1.51	1.46
3	C	803	MR0	C7-N7A	3.35	1.39	1.33
3	C	803	MR0	C23-C24	3.30	1.44	1.38
3	A	801	MR0	C5-C4	3.29	1.44	1.39
3	B	802	MR0	C25-C24	3.21	1.44	1.39
3	B	802	MR0	C22-C21	3.20	1.44	1.38
3	C	803	MR0	C27-C26	3.20	1.43	1.37
3	B	802	MR0	C21-C20	3.17	1.45	1.38
3	B	802	MR0	C5-C4	3.15	1.44	1.39
3	B	802	MR0	C7-N7A	3.08	1.38	1.33
3	A	801	MR0	C3-C2	3.05	1.44	1.39
3	B	802	MR0	C27-C26	3.00	1.42	1.37
3	C	803	MR0	C3-C2	2.99	1.44	1.39
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	C	803	MR0	C25-C24	2.85	1.43	1.39
3	B	802	MR0	C17-C28	2.85	1.58	1.51
3	C	803	MR0	C27-C28	2.84	1.44	1.39
3	A	801	MR0	C21-C20	2.84	1.44	1.38
3	C	803	MR0	C2-C7	-2.76	1.46	1.50
3	C	803	MR0	C8-N1	2.75	1.40	1.34
3	A	801	MR0	C27-C28	2.74	1.44	1.39
3	A	801	MR0	C10-N	2.67	1.54	1.47
3	A	801	MR0	C1-C	2.62	1.43	1.39
3	A	801	MR0	C25-C24	2.62	1.43	1.39
3	B	802	MR0	C1-C	2.58	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	MR0	C8-N1	2.57	1.40	1.34
3	B	802	MR0	C10-N	2.53	1.53	1.47
3	B	802	MR0	C23-C24	2.48	1.43	1.38
2	A	702	PO4	P-O3	-2.41	1.47	1.54
3	B	802	MR0	F26-C26	-2.34	1.30	1.36
2	C	701	PO4	P-O2	-2.34	1.47	1.54
3	B	802	MR0	C19-C20	2.33	1.56	1.51
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	C21-C20	2.30	1.43	1.38
3	A	801	MR0	C17-C28	2.24	1.56	1.51
3	C	803	MR0	C17-C28	2.24	1.56	1.51
2	A	702	PO4	P-O2	-2.18	1.48	1.54
3	A	801	MR0	C18-N2	2.17	1.50	1.47
3	C	803	MR0	O-C6	2.17	1.26	1.22
2	C	701	PO4	P-O3	-2.16	1.48	1.54
3	C	803	MR0	O7B-C7	2.15	1.28	1.24
3	A	801	MR0	C19-C20	2.13	1.56	1.51
3	A	801	MR0	O7B-C7	2.12	1.28	1.24
3	B	802	MR0	C18-N2	2.10	1.50	1.47
3	B	802	MR0	O-C6	2.04	1.26	1.22
3	C	803	MR0	F30-C30	-2.04	1.31	1.36
2	C	701	PO4	P-O4	-2.04	1.48	1.54
3	C	803	MR0	C19-C20	2.03	1.56	1.51

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	MR0	C-C6-N	5.95	126.09	118.66
3	B	802	MR0	C-C6-N	5.74	125.83	118.66
3	C	803	MR0	C-C6-N	4.89	124.77	118.66
3	C	803	MR0	C16-C15-N1	4.09	117.64	109.97
3	C	803	MR0	C17-C15-N1	-4.08	104.28	110.08
3	B	802	MR0	C31-C30-C29	-4.08	118.55	123.50
3	B	802	MR0	CO-O3-C24	4.03	126.14	117.50
3	A	801	MR0	C16-C15-N1	3.96	117.39	109.97
3	C	803	MR0	C28-C27-C26	3.89	122.14	118.75
3	B	802	MR0	C17-C15-N1	-3.75	104.76	110.08
3	A	801	MR0	C17-C15-N1	-3.74	104.76	110.08
3	B	802	MR0	C28-C27-C26	3.71	121.98	118.75
3	A	801	MR0	C28-C27-C26	3.59	121.88	118.75
3	C	803	MR0	C31-C26-C27	-3.45	119.30	123.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	803	MR0	C31-C30-C29	-3.25	119.55	123.50
3	A	801	MR0	C31-C26-C27	-3.23	119.57	123.50
3	A	801	MR0	C31-C30-C29	-3.18	119.64	123.50
3	B	802	MR0	C16-C18-N2	-3.17	107.79	111.95
3	C	803	MR0	C30-C31-C26	3.14	120.66	116.08
3	C	803	MR0	C24-C25-C20	3.07	123.11	119.76
3	B	802	MR0	C30-C31-C26	3.03	120.49	116.08
3	A	801	MR0	C30-C31-C26	3.00	120.45	116.08
3	B	802	MR0	C16-C15-N1	2.98	115.56	109.97
3	B	802	MR0	C20-C19-N2	-2.94	103.53	112.79
3	A	801	MR0	O-C6-N	-2.86	117.85	122.35
3	B	802	MR0	C23-C24-C25	-2.76	116.80	120.50
3	C	803	MR0	C23-C24-C25	-2.71	116.85	120.50
3	B	802	MR0	O-C6-C	-2.65	115.04	120.29
3	A	801	MR0	C24-C25-C20	2.63	122.64	119.76
3	A	801	MR0	CO-O3-C24	2.60	123.08	117.50
3	A	801	MR0	C20-C19-N2	-2.53	104.83	112.79
3	B	802	MR0	C31-C26-C27	-2.48	120.49	123.50
3	B	802	MR0	C22-C23-C24	2.39	122.59	118.98
3	B	802	MR0	O3-C24-C25	2.37	131.30	120.00
3	C	803	MR0	C20-C19-N2	-2.37	105.34	112.79
3	B	802	MR0	C24-C25-C20	2.34	122.31	119.76
3	A	801	MR0	C23-C24-C25	-2.32	117.38	120.50
3	C	803	MR0	CO-O3-C24	2.30	122.44	117.50
3	C	803	MR0	C16-C18-N2	-2.28	108.96	111.95
3	C	803	MR0	O-C6-N	-2.21	118.88	122.35
3	C	803	MR0	C3-C2-C1	2.14	122.16	119.65
3	A	801	MR0	O-C6-C	-2.13	116.06	120.29
3	B	802	MR0	O-C6-N	-2.04	119.13	122.35
3	A	801	MR0	C13-C10-N	2.04	120.92	112.37

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	801	MR0	C-C6-N-C9
3	A	801	MR0	O-C6-N-C10
3	A	801	MR0	O-C6-N-C9
3	A	801	MR0	C13-C10-N-C6
3	C	803	MR0	C-C6-N-C9
3	C	803	MR0	O-C6-N-C9
3	C	803	MR0	C13-C10-N-C6

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Mol	Chain	Res	Type	Atoms
3	A	801	MR0	C12-C11-C9-N
3	C	803	MR0	C12-C11-C9-N
3	A	801	MR0	C-C6-N-C10
3	B	802	MR0	C-C6-N-C9
3	B	802	MR0	C12-C11-C9-N
3	B	802	MR0	O-C6-N-C9
3	B	802	MR0	C13-C10-N-C6
3	C	803	MR0	C3-C2-C7-N7A
3	C	803	MR0	C3-C2-C7-O7B
3	C	803	MR0	O-C6-N-C10
3	C	803	MR0	C1-C2-C7-O7B
3	C	803	MR0	C1-C2-C7-N7A
3	A	801	MR0	C13-C10-N-C9
3	A	801	MR0	C11-C9-N-C10
3	C	803	MR0	C13-C10-N-C9
3	C	803	MR0	C-C6-N-C10
3	B	802	MR0	O-C6-N-C10
3	A	801	MR0	C11-C9-N-C6
3	C	803	MR0	C11-C9-N-C10
3	B	802	MR0	C13-C10-N-C9
3	B	802	MR0	C-C6-N-C10
3	C	803	MR0	C15-C17-C28-C27
3	C	803	MR0	C15-C17-C28-C29
3	B	802	MR0	C15-C17-C28-C27
3	B	802	MR0	C15-C17-C28-C29
3	C	803	MR0	C11-C9-N-C6

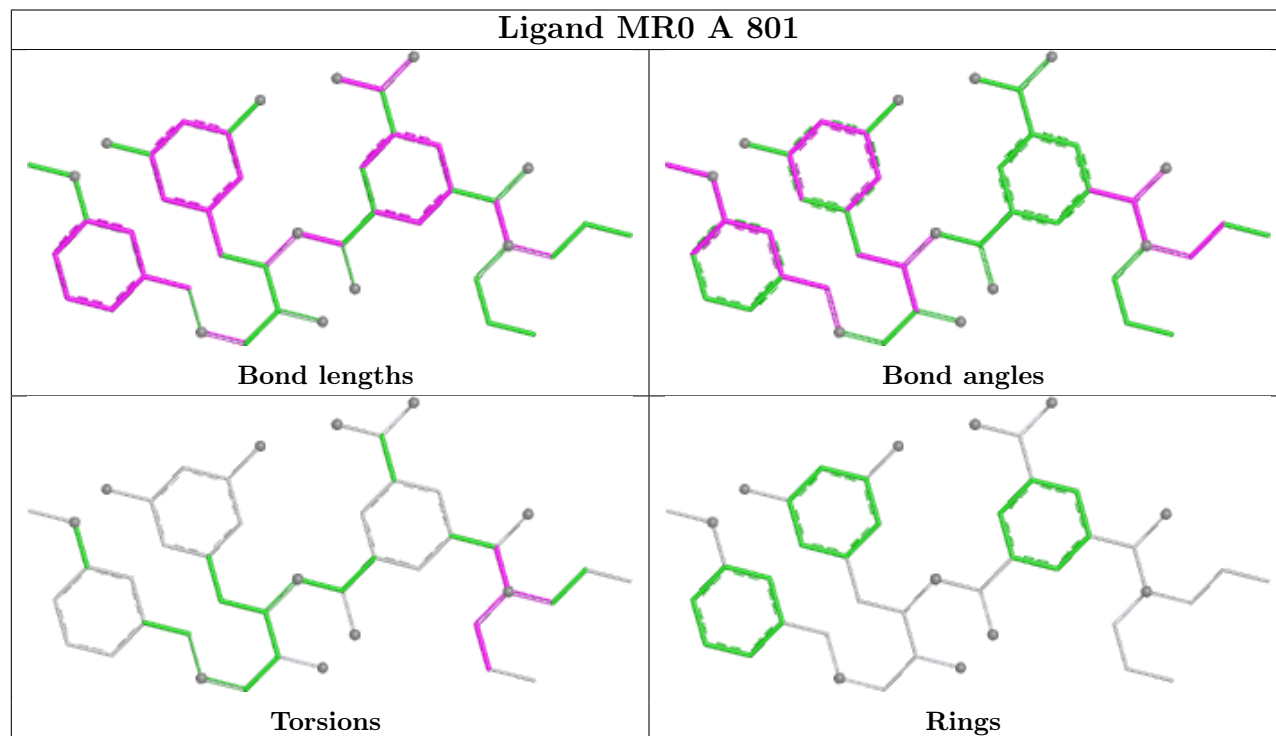
There are no ring outliers.

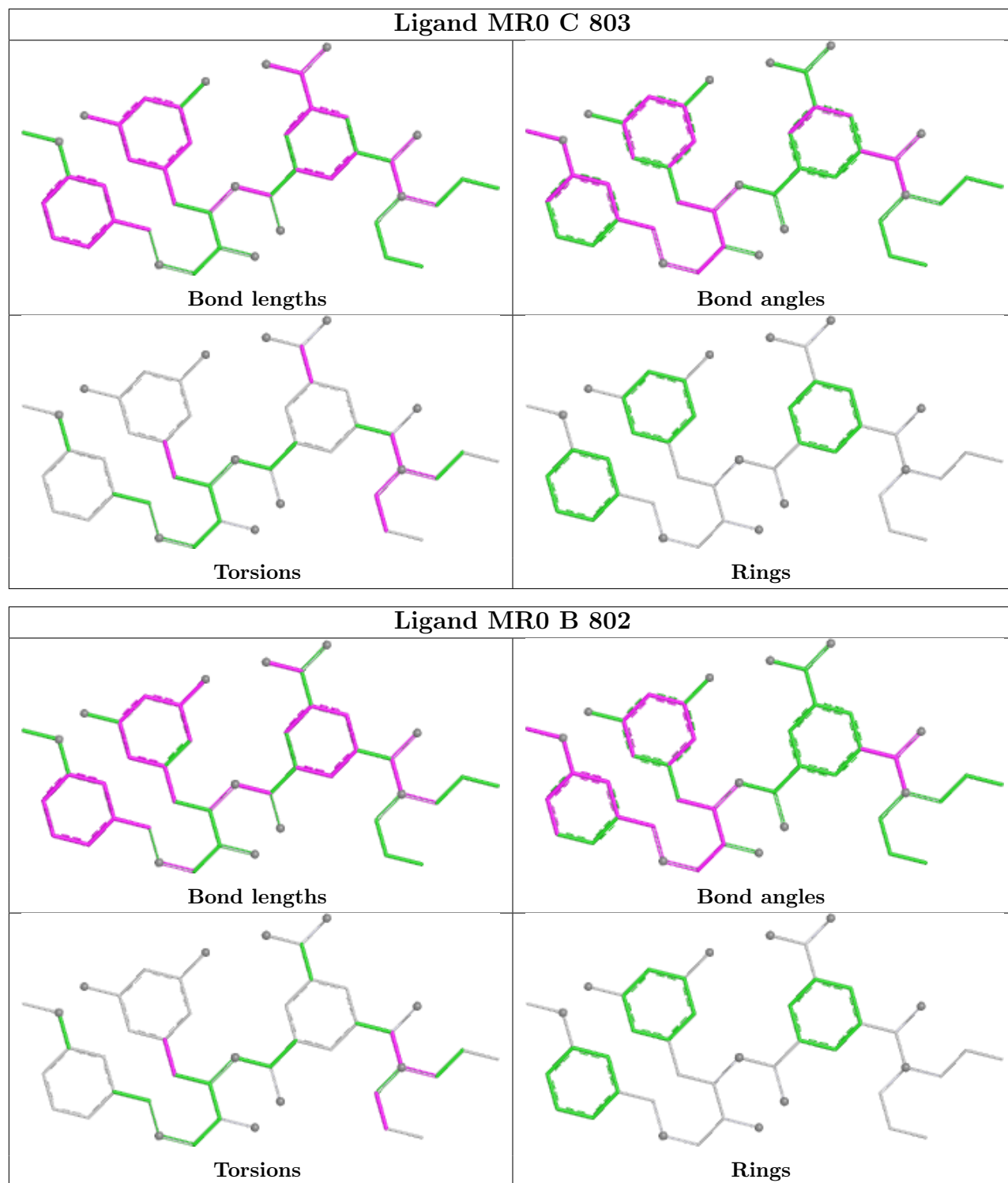
6 monomers are involved in 11 short contacts:

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

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	371/455 (81%)	-0.21	6 (1%) 70 67	26, 41, 65, 81	0
1	B	372/455 (81%)	-0.04	13 (3%) 47 42	23, 46, 74, 89	0
1	C	373/455 (81%)	-0.14	11 (2%) 53 49	26, 44, 70, 86	0
All	All	1116/1365 (81%)	-0.13	30 (2%) 56 51	23, 44, 71, 89	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	62(P)	PHE	5.1
1	B	61(P)	SER	4.8
1	A	63(P)	VAL	4.5
1	A	62(P)	PHE	4.4
1	C	61(P)	SER	3.9
1	B	62(P)	PHE	3.9
1	A	61(P)	SER	3.4
1	A	365	PHE	3.3
1	B	361	VAL	3.3
1	B	168	ALA	3.1
1	C	365	PHE	2.8
1	A	170	VAL	2.7
1	C	68	TYR	2.7
1	C	168	ALA	2.5
1	C	362	HIS	2.5
1	C	364	GLU	2.4
1	B	386	ILE	2.4
1	B	364	GLU	2.4
1	B	257	PHE	2.3
1	A	362	HIS	2.2
1	C	272	ALA	2.2
1	C	361	VAL	2.1
1	B	362	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	254	THR	2.1
1	C	254	THR	2.1
1	B	310	GLU	2.1
1	B	263	LEU	2.1
1	C	268	VAL	2.0
1	B	251	ALA	2.0
1	B	49	HIS	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](#)

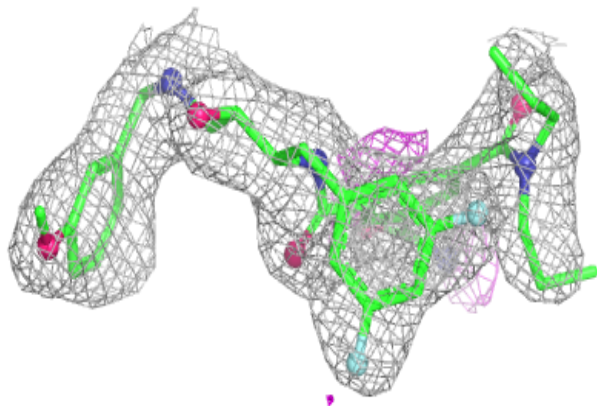
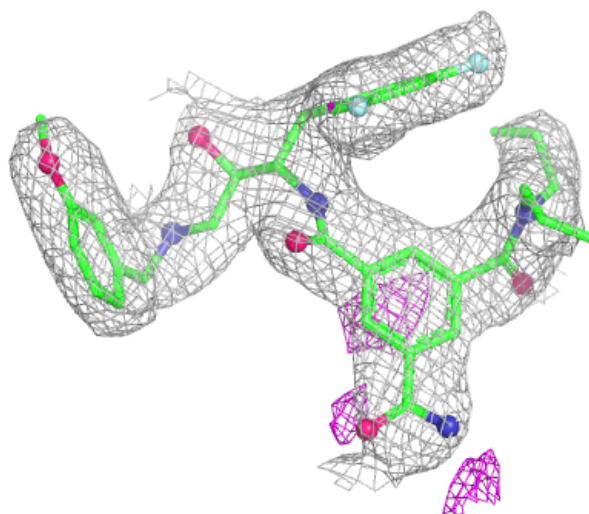
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	B	704	5/5	0.45	0.23	99,99,99,99	0
2	PO4	C	701	5/5	0.61	0.20	99,99,99,99	0
2	PO4	A	703	5/5	0.66	0.18	96,96,97,98	0
2	PO4	A	702	5/5	0.73	0.20	99,99,99,99	0
3	MR0	B	802	44/44	0.91	0.12	38,43,45,49	0
3	MR0	A	801	44/44	0.94	0.09	29,35,39,43	0
3	MR0	C	803	44/44	0.94	0.10	36,40,45,49	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.

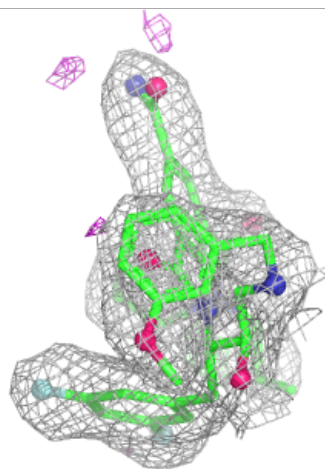
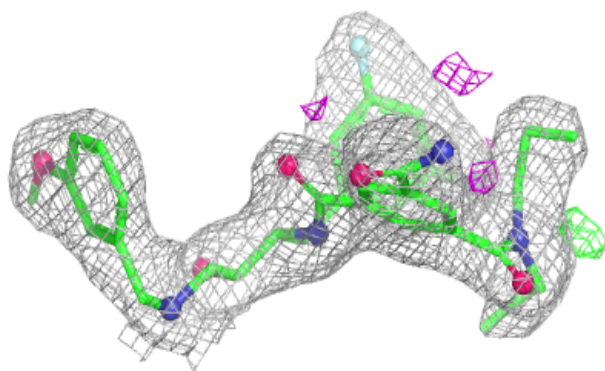
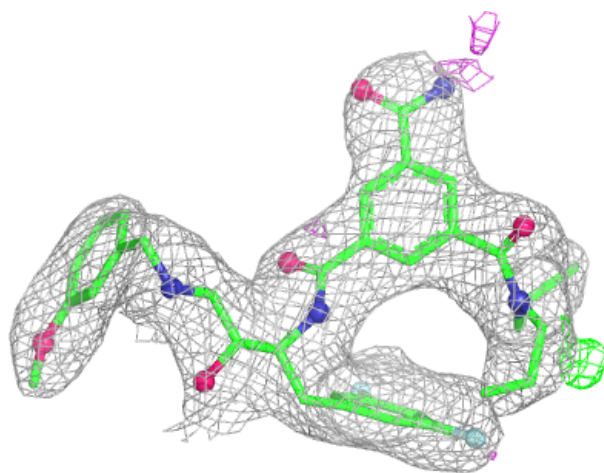
Electron density around MR0 B 802:

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



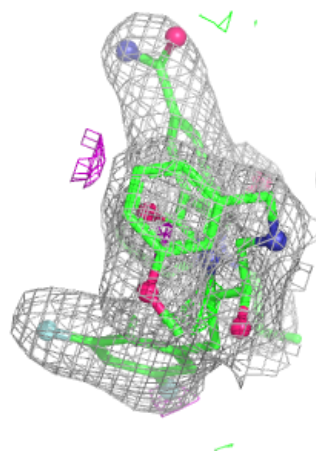
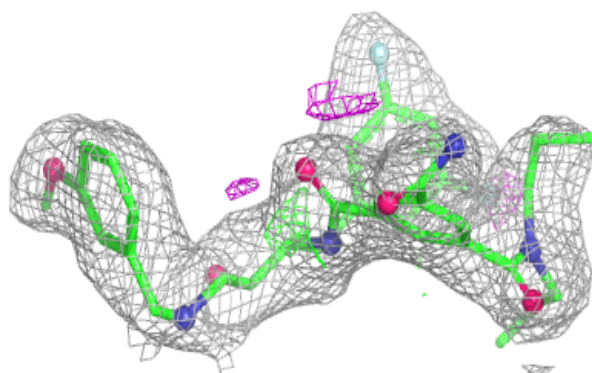
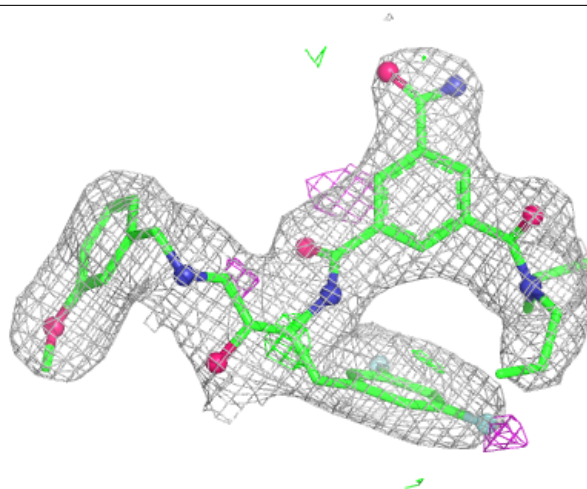
Electron density around MR0 A 801:

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



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)



6.5 Other polymers [i](#)

There are no such residues in this entry.