



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:19 PM UTC

PDB ID : 3MW0 / pdb_00003mw0
Title : X-ray structure of the doubly hydroxylated iron complex-NikA species, NikA1/O2
Authors : Cavazza, C.; Bochot, C.; Rousselot-Pailley, P.; Carpentier, P.; Cherrier, M.V.; Martin, L.; Marchi-Delapierre, C.; Fontecilla-Camps, J.C.; Menage, S.
Deposited on : 2010-05-05
Resolution : 2.30 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

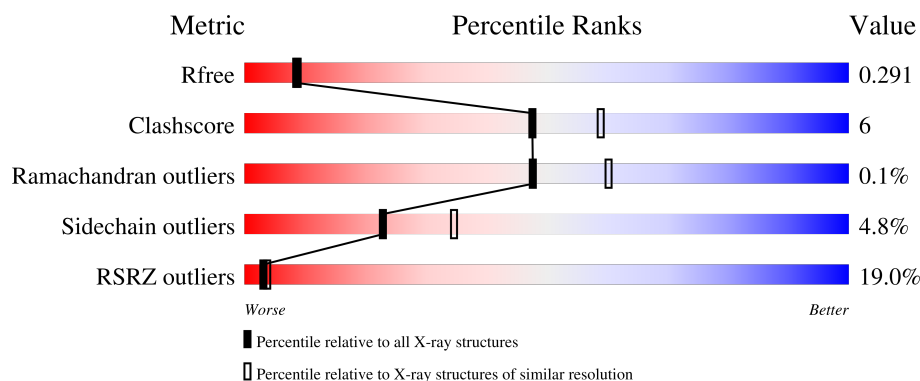
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	502	15% 86% 11% ..
1	B	502	22% 85% 13% ..

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8443 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.

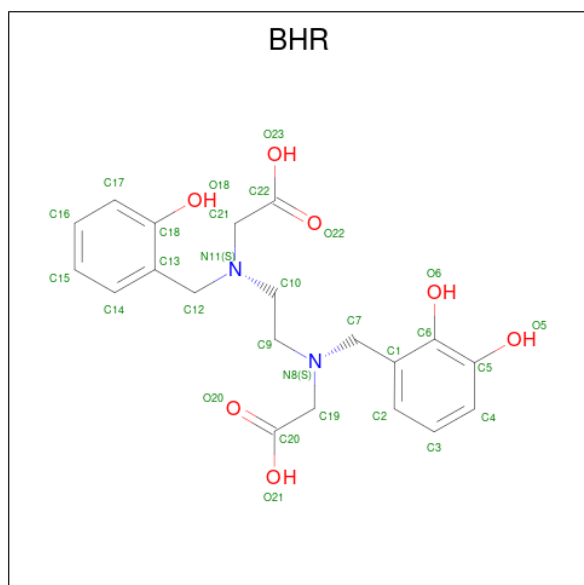
- Molecule 1 is a protein called Nickel-binding periplasmic protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	498	Total	C	N	O	S	6	2	0
			3957	2537	668	742	10			
1	B	497	Total	C	N	O	S	5	1	0
			3933	2521	666	736	10			

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

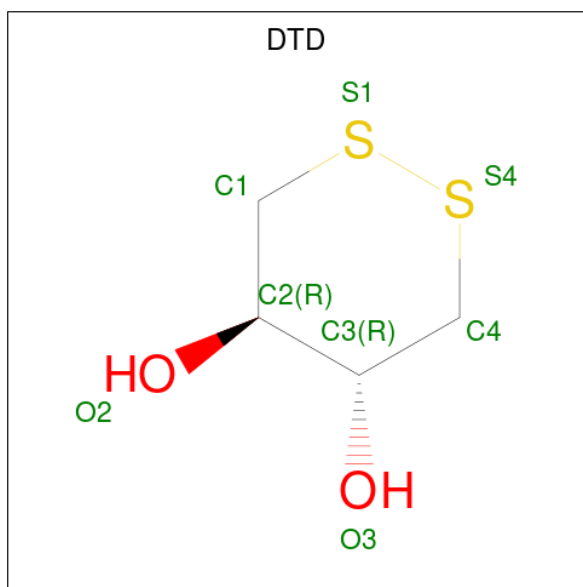
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		
2	B	1	Total	Fe	0	0
			1	1		

- Molecule 3 is 2-[2-[carboxymethyl-[(2-hydroxyphenyl)methyl]amino]ethyl-[(2,3-dihydroxyphenyl)methyl]amino]ethanoic acid (CCD ID: BHR) (formula: C₂₀H₂₄N₂O₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			29	20	2	7		
3	B	1	Total	C	N	O	0	0
			29	20	2	7		

- Molecule 4 is DITHIANE DIOL (CCD ID: DTD) (formula: $C_4H_8O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			8	4	2	2		
4	B	1	Total	C	O	S	0	0
			8	4	2	2		

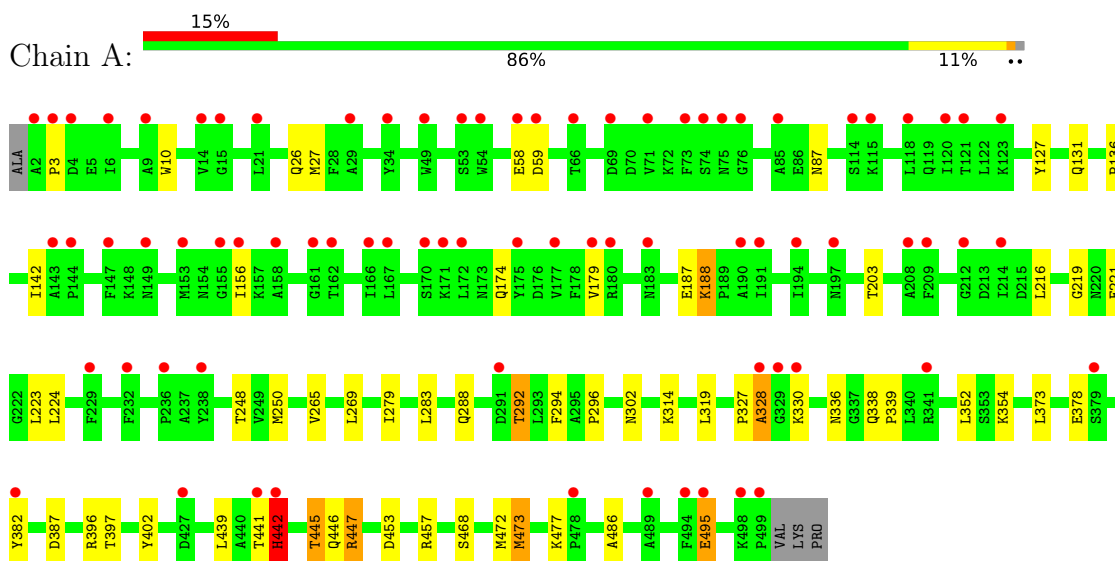
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	273	Total	O	0	0
			273	273		
5	B	204	Total	O	0	0
			204	204		

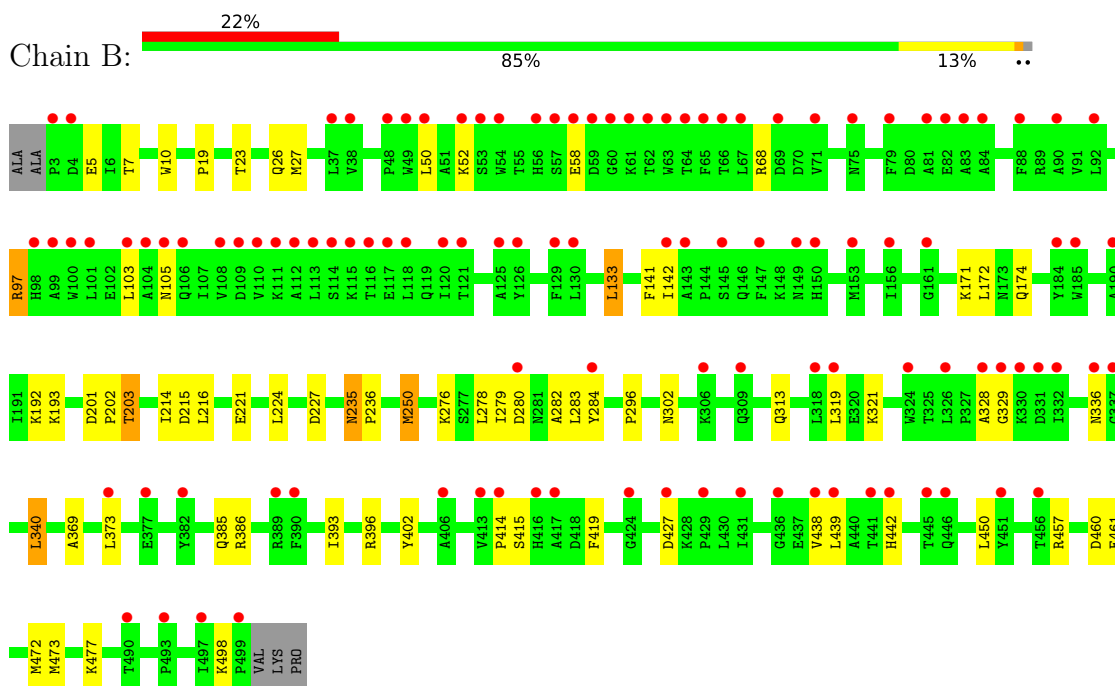
3 Residue-property plots [i](#)

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: Nickel-binding periplasmic protein



- Molecule 1: Nickel-binding periplasmic protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.98Å 94.51Å 125.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.26 – 2.30 47.26 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.26-2.30) 100.0 (47.26-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.03 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.174 , 0.244 (Not available) , 0.291	Depositor DCC
R_{free} test set	2377 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	24.7	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 28.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	8443	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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: FE, BHR, DTD

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	0.73	0/4068	0.86	5/5544 (0.1%)
1	B	0.68	0/4039	0.86	4/5502 (0.1%)
All	All	0.71	0/8107	0.86	9/11046 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	442[A]	HIS	CA-C-O	6.47	126.13	118.69
1	A	442[B]	HIS	CA-C-O	6.47	126.13	118.69
1	A	330	LYS	N-CA-C	5.50	118.95	110.42
1	B	386	ARG	CB-CG-CD	-5.30	99.10	111.30
1	B	141	PHE	N-CA-C	5.18	117.08	108.02
1	A	486	ALA	CA-C-N	5.07	125.09	119.32
1	A	486	ALA	C-N-CA	5.07	125.09	119.32
1	B	235	ASN	CA-C-N	5.05	124.77	119.82
1	B	235	ASN	C-N-CA	5.05	124.77	119.82

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	495[B]	GLU	Mainchain

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	3957	0	3898	49	1
1	B	3933	0	3862	38	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	29	0	20	0	0
3	B	29	0	20	1	0
4	A	8	0	8	1	0
4	B	8	0	8	0	0
5	A	273	0	0	7	2
5	B	204	0	0	4	0
All	All	8443	0	7816	88	3

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 6.

All (88) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:GLU:CB	1:A:59:ASP:O	1.66	1.44
1:A:292:THR:CG2	1:A:294:PHE:O	2.20	0.89
1:B:105:ASN:HB3	5:B:664:HOH:O	1.76	0.85
3:B:503:BHR:H15	5:B:551:HOH:O	1.75	0.83
1:A:58:GLU:HB2	1:A:59:ASP:O	0.76	0.81
1:A:10:TRP:HE1	1:A:26:GLN:HE21	1.31	0.75
1:A:58:GLU:CG	1:A:59:ASP:O	2.34	0.74
1:A:445:THR:HG21	5:B:618:HOH:O	1.87	0.72
1:B:438:VAL:HG12	1:B:450:LEU:HB2	1.72	0.72
1:A:188:LYS:HE3	1:A:188:LYS:H	1.55	0.72
1:B:227:ASP:HB3	1:B:284:TYR:CE2	2.26	0.70
1:A:58:GLU:CD	1:A:59:ASP:O	2.37	0.68
1:A:250:MET:HE1	1:A:397:THR:HG23	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ASN:HD21	1:A:142:ILE:H	1.43	0.64
1:A:292:THR:HG21	1:A:294:PHE:O	1.96	0.63
1:B:250:MET:HE3	1:B:396:ARG:HA	1.80	0.63
1:A:446:GLN:HE21	1:A:446:GLN:HA	1.65	0.61
1:A:453:ASP:O	1:A:457:ARG:HG3	2.00	0.61
1:B:250:MET:CE	1:B:396:ARG:HA	2.31	0.60
1:A:131:GLN:HG2	5:A:606:HOH:O	2.02	0.60
1:B:438:VAL:HG12	1:B:450:LEU:CB	2.31	0.60
1:A:219:GLY:HA3	4:A:505:DTD:H12	1.83	0.59
1:A:10:TRP:HE1	1:A:26:GLN:NE2	2.01	0.59
1:A:327:PRO:O	1:A:328:ALA:C	2.47	0.58
1:B:10:TRP:HE1	1:B:26:GLN:HE21	1.52	0.57
1:A:292:THR:HG22	1:A:294:PHE:O	2.03	0.57
1:A:453:ASP:OD2	1:A:457:ARG:HD2	2.05	0.57
1:B:414:PRO:O	1:B:415:SER:HB2	2.04	0.57
1:A:27:MET:HE1	1:A:402:TYR:OH	2.05	0.56
1:B:215:ASP:OD1	1:B:477:LYS:HE2	2.05	0.56
1:B:442:HIS:H	1:B:442:HIS:CD2	2.21	0.56
1:B:171:LYS:HB3	1:B:174:GLN:HB2	1.88	0.56
1:B:203:THR:CG2	5:B:606:HOH:O	2.54	0.55
1:B:457:ARG:O	1:B:461:GLU:HG3	2.06	0.55
1:A:248:THR:HB	1:A:250:MET:HE1	1.89	0.55
1:B:438:VAL:CG1	1:B:450:LEU:HB3	2.37	0.55
1:B:296:PRO:HB3	1:B:302:ASN:HD22	1.72	0.55
1:B:52:LYS:HE3	1:B:68:ARG:HG2	1.88	0.54
1:A:221:GLU:OE2	1:A:396:ARG:NH2	2.30	0.54
1:A:248:THR:HG22	1:A:250:MET:HE3	1.89	0.54
1:B:27:MET:HE1	1:B:402:TYR:OH	2.09	0.53
1:A:27:MET:HG3	5:A:777:HOH:O	2.09	0.52
1:A:495[B]:GLU:CD	1:A:495[B]:GLU:H	2.16	0.52
1:B:7:THR:HG22	1:B:214:ILE:HG22	1.92	0.52
1:A:327:PRO:O	1:A:328:ALA:O	2.28	0.52
1:B:103:LEU:HD21	1:B:133:LEU:HD13	1.92	0.51
1:B:414:PRO:HA	1:B:419:PHE:CD1	2.45	0.51
1:B:340:LEU:HD13	1:B:369:ALA:HB2	1.93	0.51
1:B:5:GLU:HG2	1:B:193:LYS:HB3	1.92	0.50
1:B:336:ASN:HD22	1:B:336:ASN:N	2.10	0.50
1:A:477:LYS:HE3	5:A:599:HOH:O	2.11	0.50
1:A:58:GLU:OE2	1:A:59:ASP:O	2.29	0.50
1:A:87:ASN:ND2	1:A:142:ILE:H	2.10	0.49
1:A:446:GLN:HA	1:A:446:GLN:NE2	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:438:VAL:CG1	1:B:450:LEU:CB	2.90	0.48
1:A:27:MET:HE2	1:A:136:PRO:HG2	1.96	0.48
1:A:265:VAL:O	1:A:269:LEU:HG	2.13	0.47
1:A:219:GLY:HA2	1:A:472:MET:HE3	1.97	0.47
1:B:221:GLU:OE1	1:B:396:ARG:NH1	2.45	0.47
1:B:10:TRP:HE1	1:B:26:GLN:NE2	2.13	0.46
1:A:288:GLN:O	1:A:468:SER:HB2	2.16	0.46
1:A:248:THR:HB	1:A:250:MET:CE	2.46	0.45
1:A:279:ILE:HG12	1:A:283:LEU:HD12	1.99	0.45
1:B:473:MET:HE2	1:B:473:MET:HB3	1.69	0.45
1:B:19:PRO:HG3	1:B:142:ILE:HB	1.97	0.45
1:A:441:THR:O	1:A:447:ARG:NH1	2.38	0.45
1:A:378:GLU:HG2	1:A:382:TYR:CE2	2.51	0.45
1:B:171:LYS:HD3	1:B:174:GLN:HE21	1.81	0.44
1:B:97:ARG:HH21	1:B:97:ARG:HB3	1.81	0.44
1:A:224:LEU:HD23	1:A:473:MET:CE	2.47	0.43
1:B:282:ALA:O	1:B:283:LEU:HD23	2.19	0.43
1:B:414:PRO:O	1:B:415:SER:CB	2.67	0.42
1:A:447:ARG:NH2	5:A:517:HOH:O	2.48	0.42
1:A:336:ASN:N	5:A:656:HOH:O	2.51	0.42
1:B:427:ASP:OD1	1:B:461:GLU:OE1	2.37	0.42
1:A:296:PRO:HB3	1:A:302:ASN:HD22	1.85	0.41
1:B:278:LEU:C	1:B:280:ASP:H	2.27	0.41
1:B:328:ALA:HA	1:B:329:GLY:HA2	1.65	0.41
1:A:219:GLY:CA	1:A:472:MET:HE3	2.51	0.41
1:A:3:PRO:HB2	5:A:736:HOH:O	2.20	0.41
1:B:201:ASP:HA	1:B:202:PRO:HD2	1.96	0.41
1:B:235:ASN:HA	1:B:236:PRO:HD2	1.88	0.41
1:A:224:LEU:HD23	1:A:473:MET:HE3	2.02	0.40
1:A:338:GLN:HA	1:A:339:PRO:HD3	1.96	0.40
1:A:445:THR:HG23	5:A:719:HOH:O	2.21	0.40
1:B:385:GLN:HG2	1:B:393:ILE:HD13	2.02	0.40
1:A:223:LEU:HB3	1:A:473:MET:HE2	2.03	0.40

All (3) 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)
5:A:700:HOH:O	5:A:770:HOH:O[3_555]	1.96	0.24
1:A:387:ASP:OD2	1:B:460:ASP:OD2[3_545]	2.09	0.11
5:A:654:HOH:O	5:A:671:HOH:O[3_545]	2.10	0.10

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	498/502 (99%)	481 (97%)	16 (3%)	1 (0%)	43	55
1	B	496/502 (99%)	480 (97%)	16 (3%)	0	100	100
All	All	994/1004 (99%)	961 (97%)	32 (3%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	ALA

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	422/425 (99%)	402 (95%)	20 (5%)	23	35
1	B	417/425 (98%)	396 (95%)	21 (5%)	22	33
All	All	839/850 (99%)	798 (95%)	41 (5%)	23	34

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	127	TYR
1	A	156	ILE
1	A	174	GLN
1	A	179	VAL
1	A	187	GLU

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Mol	Chain	Res	Type
1	A	188	LYS
1	A	203	THR
1	A	216	LEU
1	A	292	THR
1	A	314	LYS
1	A	319	LEU
1	A	352	LEU
1	A	354	LYS
1	A	373	LEU
1	A	439	LEU
1	A	442[A]	HIS
1	A	442[B]	HIS
1	A	445	THR
1	A	447	ARG
1	A	473	MET
1	B	23	THR
1	B	50	LEU
1	B	58	GLU
1	B	97	ARG
1	B	133	LEU
1	B	172	LEU
1	B	192	LYS
1	B	203	THR
1	B	216	LEU
1	B	224	LEU
1	B	250	MET
1	B	276	LYS
1	B	279	ILE
1	B	313	GLN
1	B	319	LEU
1	B	321	LYS
1	B	340	LEU
1	B	373	LEU
1	B	439	LEU
1	B	472	MET
1	B	498	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	87	ASN

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Mol	Chain	Res	Type
1	A	131	GLN
1	A	168	GLN
1	A	174	GLN
1	A	197	ASN
1	A	302	ASN
1	A	336	ASN
1	A	385	GLN
1	A	416	HIS
1	A	446	GLN
1	B	25	ASN
1	B	26	GLN
1	B	106	GLN
1	B	154	ASN
1	B	168	GLN
1	B	174	GLN
1	B	244	GLN
1	B	302	ASN
1	B	313	GLN
1	B	336	ASN
1	B	442	HIS
1	B	496	GLN

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](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BHR	B	503	2	30,30,30	0.93	0	40,40,40	1.55	5 (12%)
3	BHR	A	504	2	30,30,30	0.86	0	40,40,40	1.39	5 (12%)
4	DTD	B	504	-	8,8,8	1.22	1 (12%)	6,10,10	0.36	0
4	DTD	A	505	-	8,8,8	0.74	0	6,10,10	1.00	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	BHR	B	503	2	-	0/21/21/21	0/2/2/2
3	BHR	A	504	2	-	1/21/21/21	0/2/2/2
4	DTD	B	504	-	-	-	0/1/1/1
4	DTD	A	505	-	-	-	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	504	DTD	C1-S1	2.03	1.84	1.82

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	503	BHR	C12-C13-C18	5.04	125.44	120.45
3	B	503	BHR	C7-C1-C6	4.76	126.92	120.31
3	A	504	BHR	C7-C1-C6	3.54	125.22	120.31
3	A	504	BHR	C7-N8-C9	-3.52	103.94	111.28
3	A	504	BHR	C7-N8-C19	-3.47	104.80	111.77
3	A	504	BHR	C12-C13-C18	3.24	123.66	120.45
3	B	503	BHR	C7-C1-C2	-2.82	114.54	120.12
3	A	504	BHR	C7-C1-C2	-2.20	115.76	120.12
3	B	503	BHR	C12-C13-C14	-2.06	116.04	120.12
3	B	503	BHR	C12-N11-C10	-2.03	107.04	111.28

There are no chirality outliers.

All (1) torsion outliers are listed below:

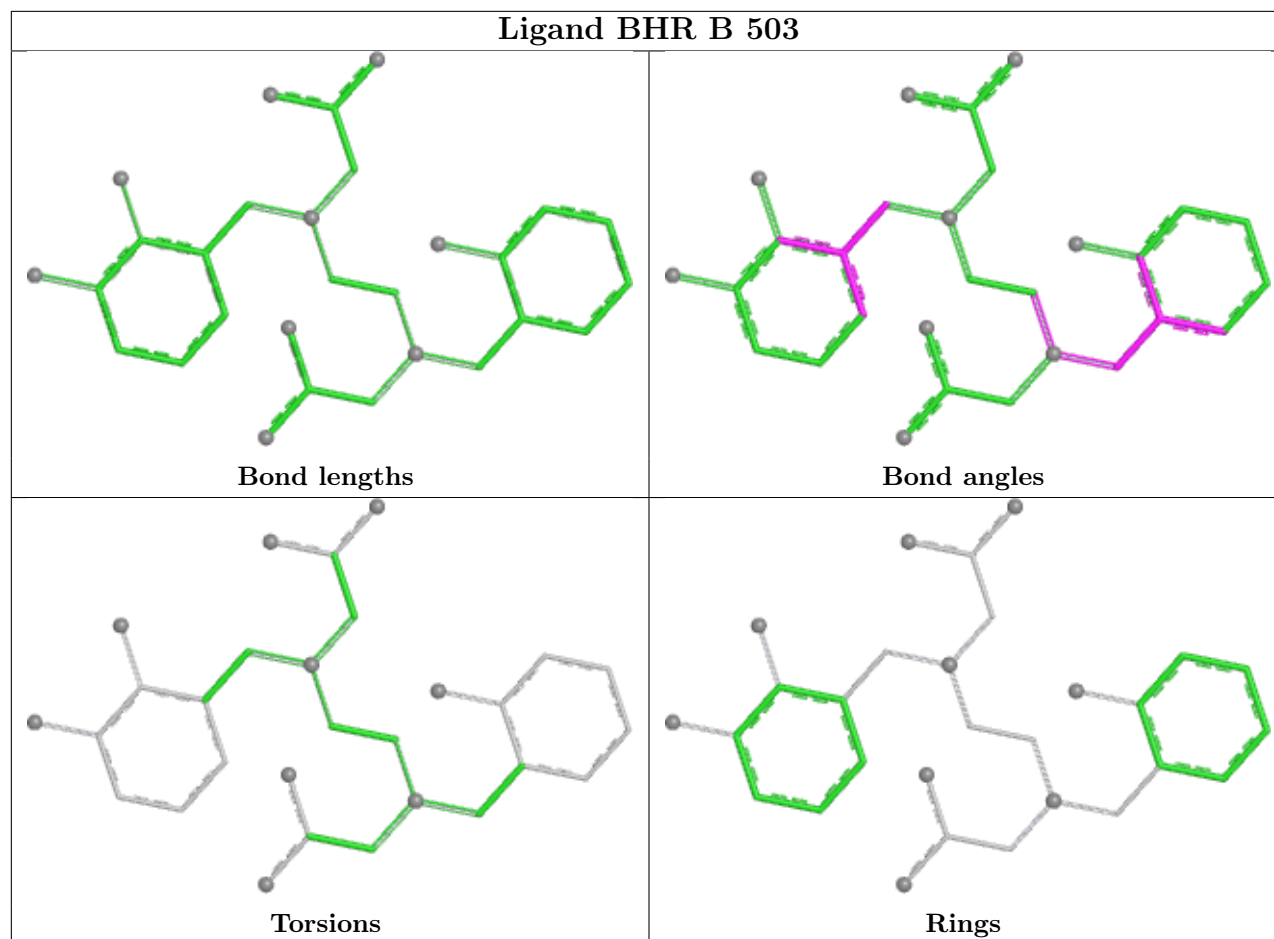
Mol	Chain	Res	Type	Atoms
3	A	504	BHR	C20-C19-N8-C9

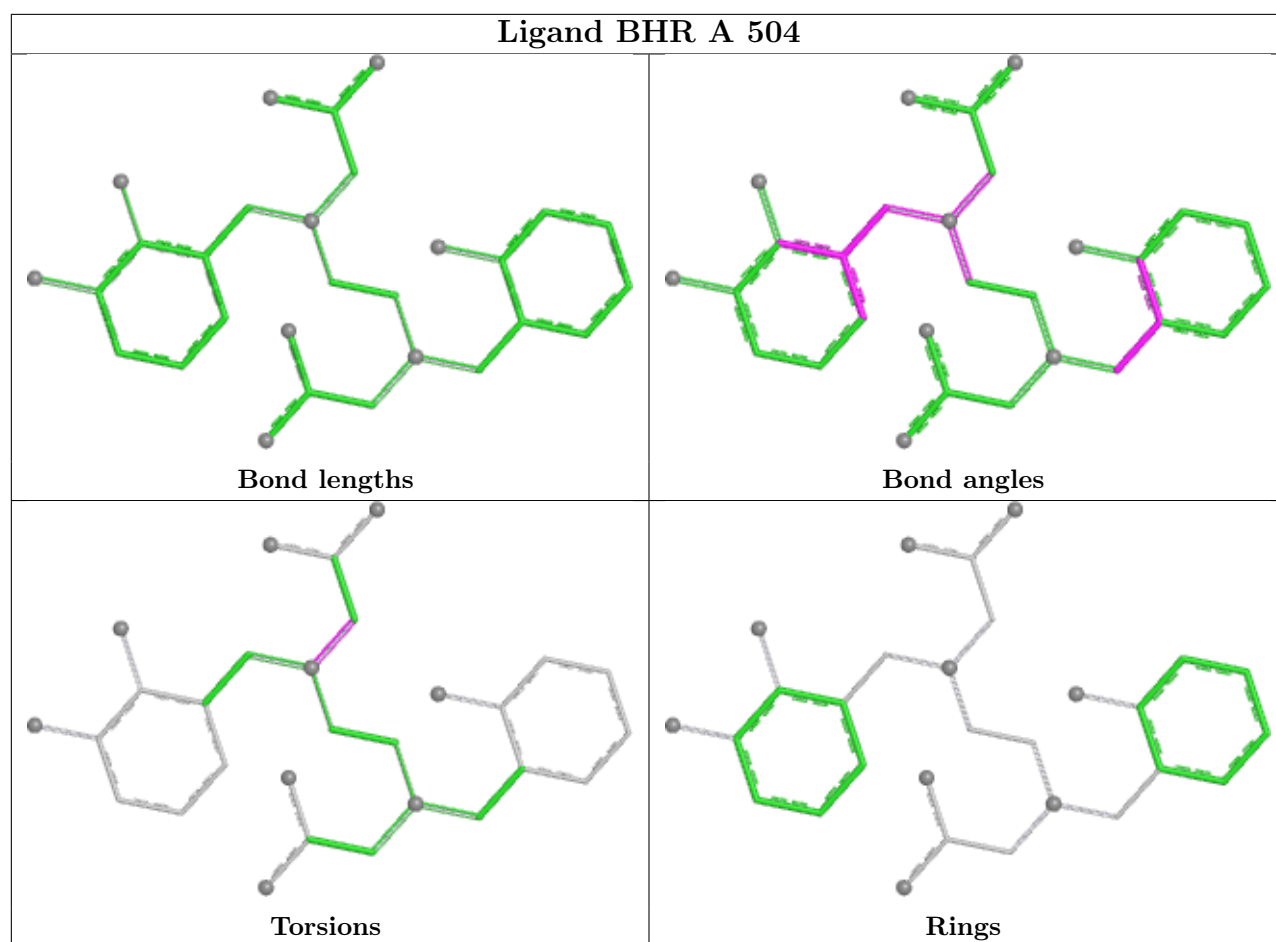
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	503	BHR	1	0
4	A	505	DTD	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	498/502 (99%)	1.06	77 (15%) 5 6	9, 19, 33, 45	4 (0%)
1	B	497/502 (99%)	1.38	112 (22%) 2 2	11, 25, 46, 60	2 (0%)
All	All	995/1004 (99%)	1.22	189 (18%) 3 4	9, 22, 42, 60	6 (0%)

All (189) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	59	ASP	6.8
1	B	442	HIS	5.1
1	A	442[A]	HIS	4.7
1	B	499	PRO	4.5
1	A	76	GLY	4.4
1	B	382	TYR	4.4
1	A	177	VAL	4.3
1	A	328	ALA	4.2
1	B	328	ALA	4.2
1	B	113	LEU	3.9
1	B	337	GLY	3.8
1	A	499	PRO	3.8
1	A	2	ALA	3.7
1	B	3	PRO	3.7
1	A	172	LEU	3.7
1	B	58	GLU	3.6
1	B	120	ILE	3.6
1	B	65	PHE	3.5
1	A	3	PRO	3.3
1	B	83	ALA	3.3
1	B	109	ASP	3.3
1	B	62	THR	3.3
1	B	121	THR	3.2
1	A	190	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	112	ALA	3.2
1	B	116	THR	3.2
1	B	329	GLY	3.2
1	A	180	ARG	3.1
1	B	52	LYS	3.1
1	A	15	GLY	3.1
1	B	330	LYS	3.1
1	B	64	THR	3.1
1	B	441	THR	3.1
1	B	145	SER	3.1
1	B	324	TRP	3.0
1	B	79	PHE	3.0
1	B	389[A]	ARG	3.0
1	B	108	VAL	3.0
1	A	291	ASP	3.0
1	B	57	SER	3.0
1	A	144	PRO	2.9
1	B	117	GLU	2.9
1	A	149	ASN	2.9
1	B	490	THR	2.9
1	A	75	ASN	2.9
1	B	336	ASN	2.9
1	A	208	ALA	2.8
1	B	63	TRP	2.8
1	A	166	ILE	2.8
1	B	105	ASN	2.8
1	B	414	PRO	2.8
1	B	318	LEU	2.8
1	B	110	VAL	2.8
1	A	494	PHE	2.8
1	B	436	GLY	2.8
1	B	118	LEU	2.8
1	B	61	LYS	2.8
1	B	280	ASP	2.7
1	B	429	PRO	2.7
1	B	92	LEU	2.7
1	A	209	PHE	2.7
1	B	71	VAL	2.7
1	B	451	TYR	2.7
1	A	212	GLY	2.7
1	B	114	SER	2.7
1	B	284	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	161	GLY	2.7
1	B	88	PHE	2.6
1	B	331	ASP	2.6
1	B	66	THR	2.6
1	B	184	TYR	2.6
1	A	379	SER	2.6
1	B	50	LEU	2.6
1	B	67	LEU	2.6
1	A	153	MET	2.6
1	A	236	PRO	2.6
1	A	175	TYR	2.6
1	A	382	TYR	2.6
1	B	56	HIS	2.6
1	A	143	ALA	2.5
1	B	390	PHE	2.5
1	A	115	LYS	2.5
1	B	150	HIS	2.5
1	A	329	GLY	2.5
1	A	170	SER	2.5
1	B	53	SER	2.5
1	B	326	LEU	2.5
1	B	90	ALA	2.5
1	A	162	THR	2.5
1	B	445	THR	2.5
1	B	129	PHE	2.5
1	B	438	VAL	2.5
1	B	37	LEU	2.5
1	B	427	ASP	2.5
1	B	81	ALA	2.5
1	B	84	ALA	2.5
1	B	125	ALA	2.5
1	B	156	ILE	2.4
1	B	497	ILE	2.4
1	A	21	LEU	2.4
1	B	59	ASP	2.4
1	B	75	ASN	2.4
1	A	158	ALA	2.4
1	A	74	SER	2.4
1	B	49	TRP	2.4
1	A	85	ALA	2.4
1	B	424	GLY	2.4
1	A	114	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	179	VAL	2.4
1	B	185	TRP	2.4
1	B	82	GLU	2.4
1	B	115	LYS	2.4
1	B	101	LEU	2.4
1	A	58	GLU	2.4
1	A	54	TRP	2.4
1	A	427	ASP	2.3
1	B	149	ASN	2.3
1	A	441	THR	2.3
1	A	120	ILE	2.3
1	B	98	HIS	2.3
1	B	377	GLU	2.3
1	A	330	LYS	2.3
1	B	190	ALA	2.3
1	A	191	ILE	2.3
1	B	413	VAL	2.3
1	B	104	ALA	2.3
1	B	332	ILE	2.3
1	A	14	VAL	2.3
1	A	73	PHE	2.3
1	A	341	ARG	2.3
1	B	446	GLN	2.3
1	B	54	TRP	2.2
1	B	106	GLN	2.2
1	A	156	ILE	2.2
1	A	194	ILE	2.2
1	A	167	LEU	2.2
1	B	103	LEU	2.2
1	A	161	GLY	2.2
1	A	29	ALA	2.2
1	B	417	ALA	2.2
1	A	6	ILE	2.2
1	B	431	ILE	2.2
1	B	306	LYS	2.2
1	B	416	HIS	2.2
1	B	143	ALA	2.2
1	A	53	SER	2.2
1	B	130	LEU	2.2
1	B	439	LEU	2.2
1	A	183	ASN	2.2
1	B	69	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	493	PRO	2.2
1	A	9	ALA	2.2
1	A	121	THR	2.2
1	B	456	THR	2.2
1	B	142	ILE	2.1
1	A	34	TYR	2.1
1	A	238	TYR	2.1
1	A	498	LYS	2.1
1	B	319	LEU	2.1
1	A	232	PHE	2.1
1	A	478	PRO	2.1
1	B	309	GLN	2.1
1	A	155	GLY	2.1
1	B	406	ALA	2.1
1	A	123	LYS	2.1
1	B	153	MET	2.1
1	A	197	ASN	2.1
1	B	126	TYR	2.1
1	A	71	VAL	2.1
1	A	49	TRP	2.1
1	B	147	PHE	2.1
1	B	60	GLY	2.1
1	A	489	ALA	2.1
1	B	99	ALA	2.1
1	A	4	ASP	2.1
1	A	69	ASP	2.1
1	B	38	VAL	2.1
1	B	100	TRP	2.1
1	A	66	THR	2.0
1	A	495[A]	GLU	2.0
1	A	214	ILE	2.0
1	A	147	PHE	2.0
1	A	229	PHE	2.0
1	A	118	LEU	2.0
1	B	4	ASP	2.0
1	B	373	LEU	2.0
1	B	48	PRO	2.0
1	A	171	LYS	2.0
1	B	111	LYS	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](#)

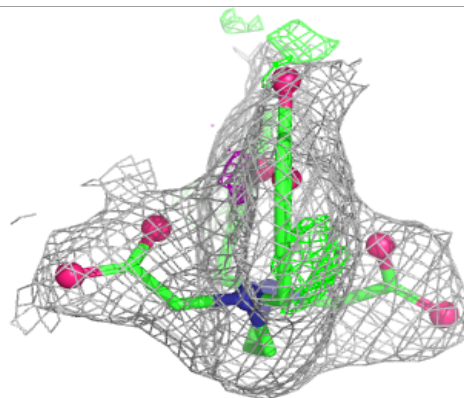
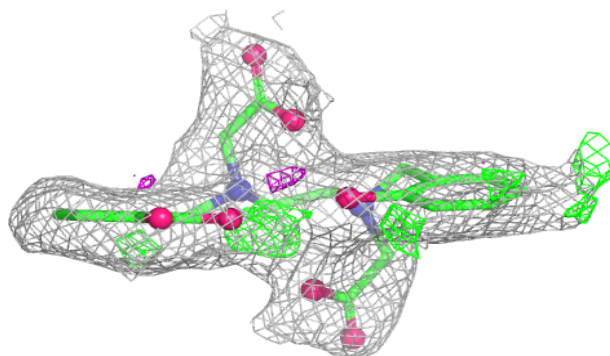
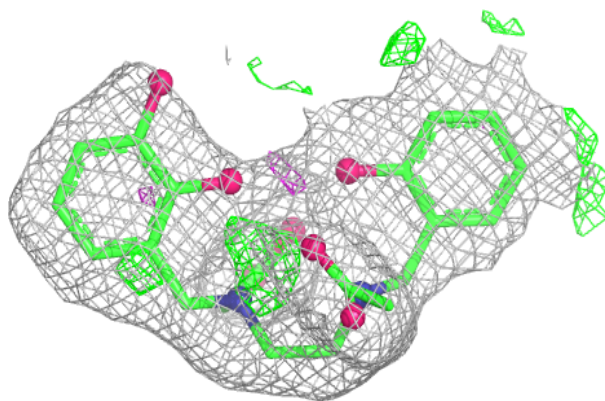
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	DTD	A	505	8/8	0.68	0.34	75,75,75,76	0
4	DTD	B	504	8/8	0.74	0.27	79,79,79,79	0
3	BHR	A	504	29/29	0.85	0.14	23,27,29,31	0
3	BHR	B	503	29/29	0.85	0.13	19,27,31,32	0
2	FE	A	503	1/1	0.92	0.06	20,20,20,20	0
2	FE	B	505	1/1	0.92	0.06	23,23,23,23	0

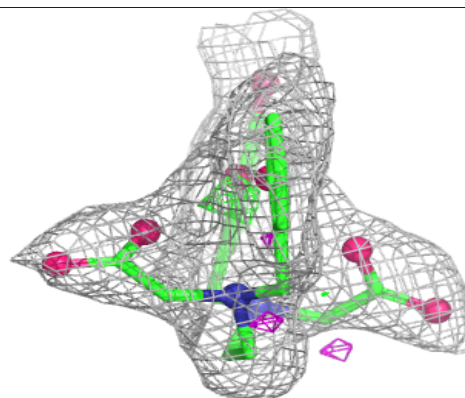
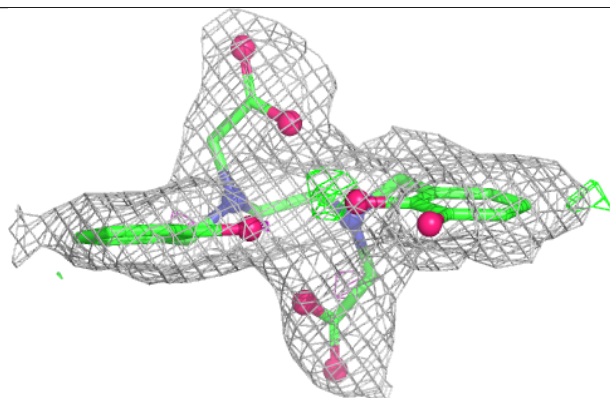
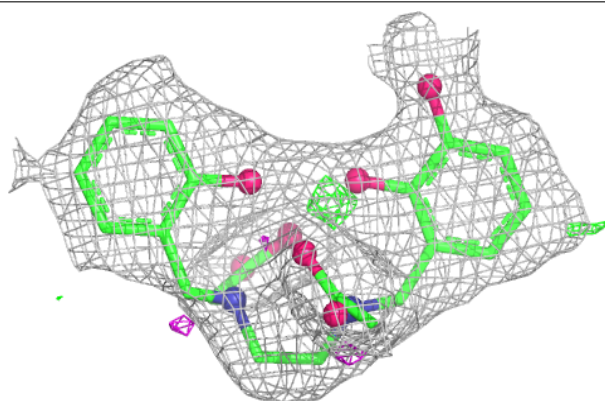
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BHR A 504:

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

**Electron density around BHR B 503:**

$2mF_o-DF_c$ (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.