



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:39 PM UTC

PDB ID : 1IYH / pdb_00001iyh
Title : Crystal structure of hematopoietic prostaglandin D synthase
Authors : Inoue, T.
Deposited on : 2002-08-26
Resolution : 1.70 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

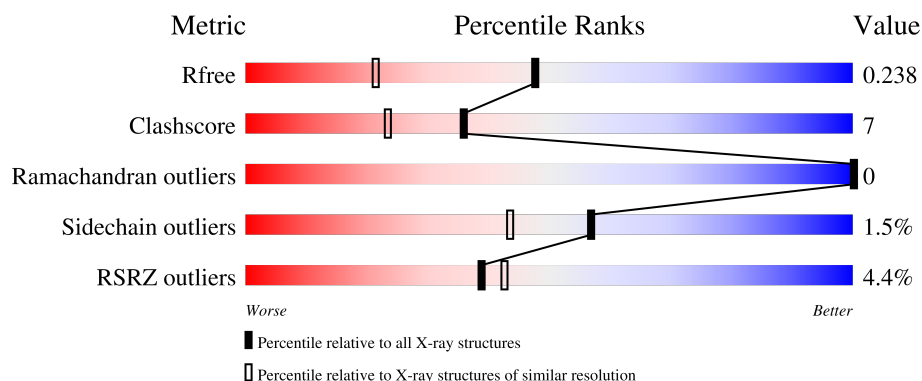
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	198	3% 83% 16% .
1	B	198	% 86% 14% .
1	C	198	5% 79% 20% .
1	D	198	9% 85% 15% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7616 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 HEMATOPOIETIC PROSTAGLANDIN D SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	0	0	0
			1638	1058	273	298	9			
1	B	198	Total	C	N	O	S	0	0	0
			1638	1058	273	298	9			
1	C	198	Total	C	N	O	S	0	0	0
			1638	1058	273	298	9			
1	D	198	Total	C	N	O	S	0	0	0
			1638	1058	273	298	9			

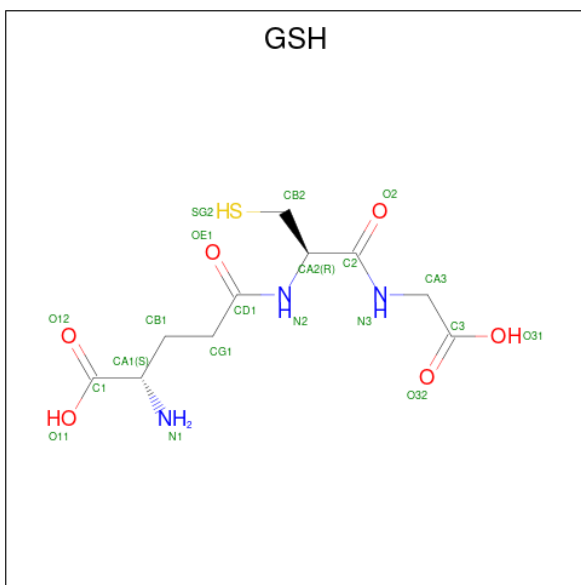
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	144	MET	ASN	conflict	UNP O60760
B	344	MET	ASN	conflict	UNP O60760
C	544	MET	ASN	conflict	UNP O60760
D	744	MET	ASN	conflict	UNP O60760

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 3 is GLUTATHIONE (CCD ID: GSH) (formula: C₁₀H₁₇N₃O₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 20	C 10	N 3	O 6	S 1	0	0
3	B	1	Total 20	C 10	N 3	O 6	S 1	0	0
3	C	1	Total 20	C 10	N 3	O 6	S 1	0	0
3	D	1	Total 20	C 10	N 3	O 6	S 1	0	0

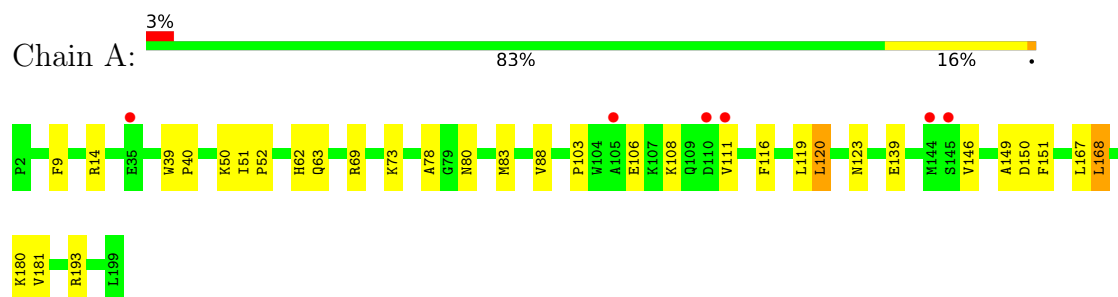
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	279	Total O 279 279	0	0
4	B	268	Total O 268 268	0	0
4	C	234	Total O 234 234	0	0
4	D	201	Total O 201 201	0	0

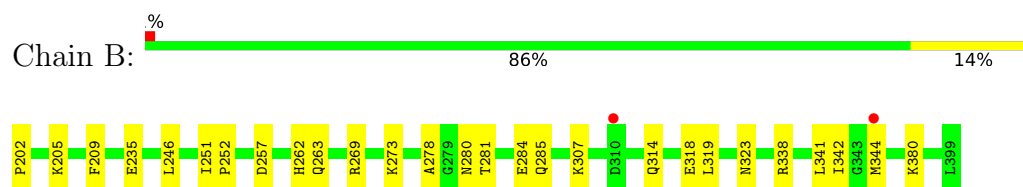
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

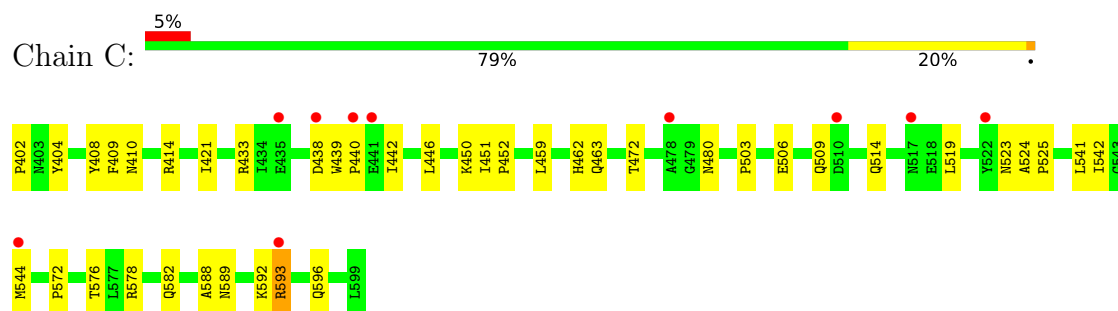
• Molecule 1: HEMATOPOIETIC PROSTAGLANDIN D SYNTHASE



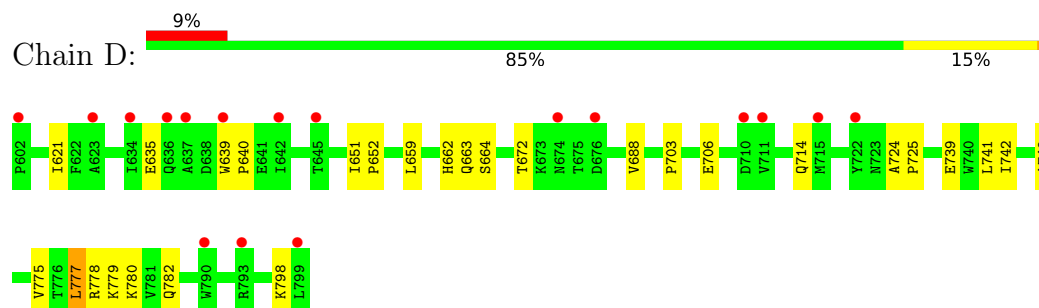
• Molecule 1: HEMATOPOIETIC PROSTAGLANDIN D SYNTHASE



• Molecule 1: HEMATOPOIETIC PROSTAGLANDIN D SYNTHASE



• Molecule 1: HEMATOPOIETIC PROSTAGLANDIN D SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.11Å 47.37Å 184.70Å 90.00° 97.44° 90.00°	Depositor
Resolution (Å)	30.94 – 1.70 30.94 – 1.70	Depositor EDS
% Data completeness (in resolution range)	86.4 (30.94-1.70) 86.5 (30.94-1.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.65 (at 1.71Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.205 , 0.238 0.206 , 0.238	Depositor DCC
R_{free} test set	8029 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	18.8	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.012 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7616	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.77 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.4792e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GSH, MG

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	0.37	0/1681	0.80	2/2283 (0.1%)
1	B	0.37	0/1681	0.81	2/2283 (0.1%)
1	C	0.36	0/1681	0.78	1/2283 (0.0%)
1	D	0.35	0/1681	0.77	1/2283 (0.0%)
All	All	0.36	0/6724	0.79	6/9132 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	278	ALA	N-CA-C	6.68	118.64	111.36
1	C	523	ASN	N-CA-C	6.33	117.98	111.14
1	A	78	ALA	N-CA-C	6.32	118.25	111.36
1	A	80	ASN	N-CA-C	5.83	118.58	111.82
1	D	664	SER	N-CA-C	5.51	118.00	111.33
1	B	280	ASN	N-CA-C	5.40	119.59	112.89

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	1638	0	1622	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1638	0	1622	19	0
1	C	1638	0	1622	33	0
1	D	1638	0	1622	19	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	20	0	15	0	0
3	B	20	0	15	0	0
3	C	20	0	15	2	0
3	D	20	0	15	0	0
4	A	279	0	0	15	0
4	B	268	0	0	4	0
4	C	234	0	0	7	0
4	D	201	0	0	3	0
All	All	7616	0	6548	96	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 7.

All (96) 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:181:VAL:HG22	4:A:1934:HOH:O	1.75	0.84
1:C:576:THR:HG23	1:D:765:PRO:HG3	1.62	0.80
1:C:593:ARG:HH11	1:C:593:ARG:HG3	1.54	0.71
1:B:319:LEU:HA	1:B:323:ASN:HD22	1.58	0.68
1:B:338:ARG:HD2	4:B:2168:HOH:O	1.93	0.68
1:A:151:PHE:HA	4:A:1957:HOH:O	1.92	0.67
1:C:442:ILE:O	1:C:446:LEU:HD13	1.93	0.67
1:A:151:PHE:HD1	4:A:1957:HOH:O	1.78	0.65
1:C:589:ASN:HA	1:C:592:LYS:NZ	2.15	0.62
1:B:281:THR:O	1:B:285:GLN:HG3	2.01	0.60
1:A:9:PHE:CE1	1:A:51:ILE:HD11	2.37	0.59
1:B:209:PHE:CE1	1:B:251:ILE:HD11	2.40	0.57
1:A:119:LEU:HA	1:A:123:ASN:HD22	1.69	0.57
1:D:703:PRO:HB2	1:D:706:GLU:HB2	1.88	0.56
1:C:402:PRO:HG2	1:C:404:TYR:CZ	2.41	0.56
1:C:524:ALA:HB3	1:C:525:PRO:HD3	1.87	0.56
1:A:116:PHE:O	1:A:120:LEU:HD22	2.06	0.56
1:C:410:ASN:HD22	1:C:433:ARG:HH11	1.53	0.55
1:A:50:LYS:HE2	4:A:2019:HOH:O	2.06	0.55
1:A:39:TRP:HB3	1:A:40:PRO:HD3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:VAL:CG2	4:A:1957:HOH:O	2.54	0.55
1:C:440:PRO:HG2	4:C:1705:HOH:O	2.07	0.55
1:C:503:PRO:HB2	1:C:506:GLU:HB2	1.89	0.55
1:C:439:TRP:N	1:C:440:PRO:HD2	2.22	0.55
1:C:593:ARG:HG3	1:C:593:ARG:NH1	2.20	0.54
1:C:480:ASN:ND2	4:C:1815:HOH:O	2.40	0.54
1:D:639:TRP:HB3	1:D:640:PRO:HD3	1.88	0.54
1:D:778:ARG:O	1:D:782:GLN:HG3	2.07	0.54
4:A:2060:HOH:O	1:C:462:HIS:HE1	1.92	0.53
1:A:193:ARG:HD3	4:A:2115:HOH:O	2.09	0.52
1:B:281:THR:OG1	1:B:284:GLU:HG3	2.09	0.52
1:D:659:LEU:HD12	4:D:1978:HOH:O	2.08	0.52
1:C:402:PRO:HG2	1:C:404:TYR:CE1	2.45	0.52
1:A:181:VAL:CG1	4:A:1957:HOH:O	2.58	0.52
1:C:589:ASN:HA	1:C:592:LYS:HZ3	1.75	0.51
1:A:150:ASP:C	4:A:1957:HOH:O	2.54	0.51
1:D:651:ILE:HB	1:D:652:PRO:HA	1.93	0.51
1:D:621:ILE:HD13	1:D:672:THR:HG23	1.94	0.50
1:D:714:GLN:HB3	4:D:1958:HOH:O	2.12	0.50
1:D:724:ALA:HB3	1:D:725:PRO:HD3	1.94	0.50
1:A:14:ARG:NH1	4:A:2132:HOH:O	2.42	0.49
1:B:205:LYS:HB2	4:B:2120:HOH:O	2.12	0.49
1:B:262:HIS:HE1	4:D:1869:HOH:O	1.95	0.49
1:B:202:PRO:HB3	1:B:257:ASP:OD1	2.12	0.49
1:B:235:GLU:OE2	1:B:235:GLU:HA	2.13	0.49
1:A:103:PRO:HB2	1:A:106:GLU:HB2	1.95	0.49
1:C:578:ARG:O	1:C:582:GLN:HG3	2.13	0.49
1:A:108:LYS:HB2	1:A:111:VAL:HG12	1.94	0.49
1:A:14:ARG:HD3	4:A:2132:HOH:O	2.13	0.48
1:A:151:PHE:N	4:A:1957:HOH:O	2.45	0.48
1:D:741:LEU:C	1:D:742:ILE:HD12	2.39	0.48
1:A:62:HIS:HE1	4:C:1652:HOH:O	1.96	0.48
1:B:341:LEU:HB3	1:B:342:ILE:HD12	1.96	0.48
1:B:319:LEU:HD23	1:B:323:ASN:HD22	1.79	0.47
1:A:62:HIS:O	1:A:63:GLN:HB2	2.13	0.47
1:C:596:GLN:HA	4:C:1730:HOH:O	2.14	0.47
1:B:319:LEU:HD23	1:B:323:ASN:ND2	2.29	0.46
1:A:116:PHE:O	1:A:120:LEU:CD2	2.64	0.46
1:A:151:PHE:CA	4:A:1957:HOH:O	2.57	0.46
1:C:451:ILE:HB	1:C:452:PRO:HA	1.97	0.46
1:A:83:MET:HE2	1:C:459:LEU:HD13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:GLN:O	1:B:318:GLU:HG3	2.16	0.45
1:C:462:HIS:O	1:C:463:GLN:HB2	2.16	0.45
1:B:380:LYS:HD3	4:B:2142:HOH:O	2.16	0.45
1:C:588:ALA:O	1:C:592:LYS:HG2	2.16	0.45
1:C:414:ARG:NH2	4:C:1768:HOH:O	2.49	0.45
1:B:341:LEU:CB	1:B:342:ILE:HD12	2.46	0.44
1:D:688:VAL:HG13	1:D:749:ALA:HB2	2.00	0.44
1:C:462:HIS:HD2	4:C:1636:HOH:O	2.00	0.44
1:A:51:ILE:HB	1:A:52:PRO:HA	1.99	0.44
1:A:88:VAL:HG13	1:A:149:ALA:HB2	1.98	0.44
1:A:139:GLU:HG2	1:A:180:LYS:HE2	1.99	0.43
1:C:572:PRO:HB3	1:D:798:LYS:HA	1.99	0.43
1:D:742:ILE:HD12	1:D:742:ILE:N	2.34	0.43
1:A:167:LEU:HG	1:A:168:LEU:HD13	2.00	0.43
1:B:307:LYS:HB2	4:B:2044:HOH:O	2.18	0.43
1:C:409:PHE:CE1	1:C:451:ILE:HD11	2.54	0.43
1:C:414:ARG:HD2	3:C:1600:GSH:SG2	2.58	0.43
1:C:421:ILE:HD13	1:C:472:THR:HG23	2.00	0.43
1:D:775:VAL:O	1:D:779:LYS:HG2	2.18	0.42
1:C:450:LYS:HB3	3:C:1600:GSH:HG13	2.02	0.42
1:C:519:LEU:HD22	4:C:1658:HOH:O	2.19	0.42
1:D:739:GLU:HG3	1:D:780:LYS:HE3	2.02	0.42
1:C:542:ILE:HD12	1:C:542:ILE:N	2.35	0.41
1:B:269:ARG:O	1:B:273:LYS:HG3	2.21	0.41
1:A:69:ARG:O	1:A:73:LYS:HG3	2.20	0.41
1:A:146:VAL:HG23	4:A:1934:HOH:O	2.21	0.41
1:B:262:HIS:O	1:B:263:GLN:HB2	2.21	0.41
1:D:753:TRP:CD1	1:D:777:LEU:HD13	2.55	0.41
1:B:251:ILE:HB	1:B:252:PRO:HA	2.02	0.41
1:C:408:TYR:CG	1:C:409:PHE:N	2.89	0.41
1:C:541:LEU:CB	1:C:542:ILE:HD12	2.51	0.41
1:D:662:HIS:O	1:D:663:GLN:HB2	2.21	0.41
1:D:741:LEU:HG	1:D:777:LEU:HD11	2.02	0.41
1:C:576:THR:CG2	1:D:765:PRO:HG3	2.42	0.40
1:A:181:VAL:HG21	4:A:1957:HOH:O	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	196/198 (99%)	191 (97%)	5 (3%)	0	100	100
1	B	196/198 (99%)	193 (98%)	3 (2%)	0	100	100
1	C	196/198 (99%)	192 (98%)	4 (2%)	0	100	100
1	D	196/198 (99%)	191 (97%)	5 (3%)	0	100	100
All	All	784/792 (99%)	767 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	178/178 (100%)	176 (99%)	2 (1%)	65	54
1	B	178/178 (100%)	176 (99%)	2 (1%)	65	54
1	C	178/178 (100%)	173 (97%)	5 (3%)	38	21
1	D	178/178 (100%)	176 (99%)	2 (1%)	65	54
All	All	712/712 (100%)	701 (98%)	11 (2%)	57	43

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	LEU
1	A	168	LEU

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Mol	Chain	Res	Type
1	B	246	LEU
1	B	344	MET
1	C	438	ASP
1	C	509	GLN
1	C	514	GLN
1	C	544	MET
1	C	593	ARG
1	D	635	GLU
1	D	777	LEU

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

Mol	Chain	Res	Type
1	A	3	ASN
1	A	36	GLN
1	A	62	HIS
1	A	109	GLN
1	A	114	GLN
1	A	123	ASN
1	A	129	GLN
1	A	170	ASN
1	A	171	HIS
1	B	262	HIS
1	B	274	ASN
1	B	323	ASN
1	B	370	ASN
1	C	403	ASN
1	C	410	ASN
1	C	462	HIS
1	C	480	ASN
1	C	509	GLN
1	C	517	ASN
1	C	523	ASN
1	C	570	ASN
1	C	596	GLN
1	D	628	GLN
1	D	662	HIS
1	D	680	ASN
1	D	709	GLN
1	D	770	ASN
1	D	796	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	GSH	A	1200	-	18,19,19	1.36	3 (16%)	21,24,24	1.14	1 (4%)
3	GSH	D	1800	-	18,19,19	1.27	2 (11%)	21,24,24	1.18	2 (9%)
3	GSH	B	1400	-	18,19,19	1.38	3 (16%)	21,24,24	1.14	2 (9%)
3	GSH	C	1600	-	18,19,19	1.27	1 (5%)	21,24,24	1.18	1 (4%)

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	GSH	A	1200	-	-	3/24/24/24	-
3	GSH	D	1800	-	-	6/24/24/24	-
3	GSH	B	1400	-	-	3/24/24/24	-
3	GSH	C	1600	-	-	5/24/24/24	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1200	GSH	O12-C1	3.70	1.33	1.22
3	D	1800	GSH	O12-C1	3.50	1.32	1.22
3	B	1400	GSH	O12-C1	3.47	1.32	1.22
3	C	1600	GSH	O12-C1	3.43	1.32	1.22
3	B	1400	GSH	CG1-CD1	2.81	1.57	1.51
3	A	1200	GSH	CG1-CD1	2.35	1.56	1.51
3	B	1400	GSH	O11-C1	-2.25	1.23	1.30
3	D	1800	GSH	O11-C1	-2.07	1.24	1.30
3	A	1200	GSH	O31-C3	-2.01	1.24	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1600	GSH	CB1-CG1-CD1	-3.74	104.72	113.06
3	D	1800	GSH	CB1-CG1-CD1	-3.69	104.84	113.06
3	A	1200	GSH	CB1-CG1-CD1	-3.49	105.28	113.06
3	B	1400	GSH	CB1-CG1-CD1	-3.32	105.66	113.06
3	B	1400	GSH	CB2-CA2-C2	-2.72	103.59	109.50
3	D	1800	GSH	CB2-CA2-C2	-2.14	104.85	109.50

There are no chirality outliers.

All (17) torsion outliers are listed below:

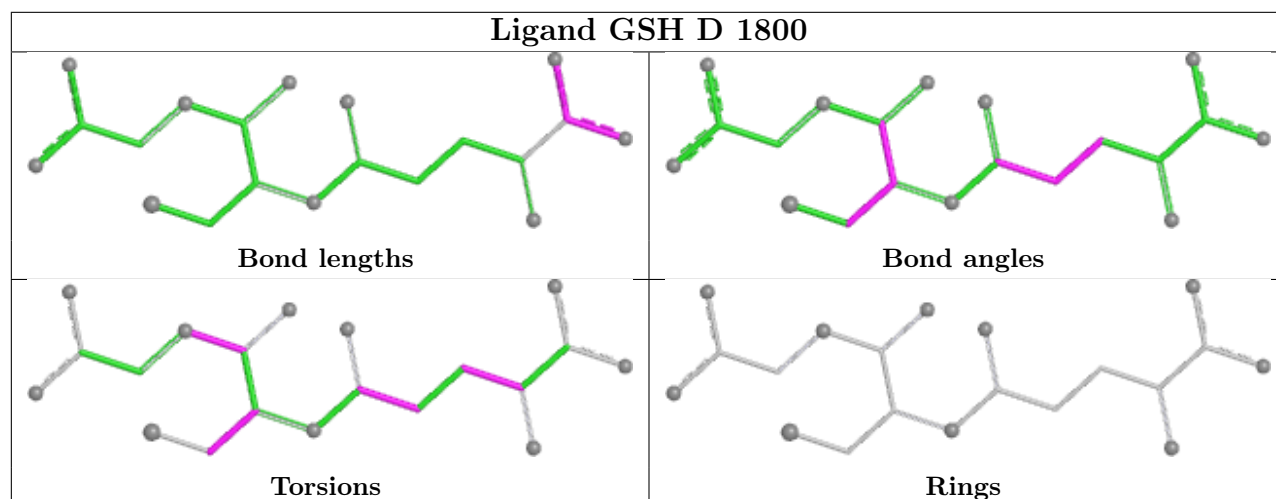
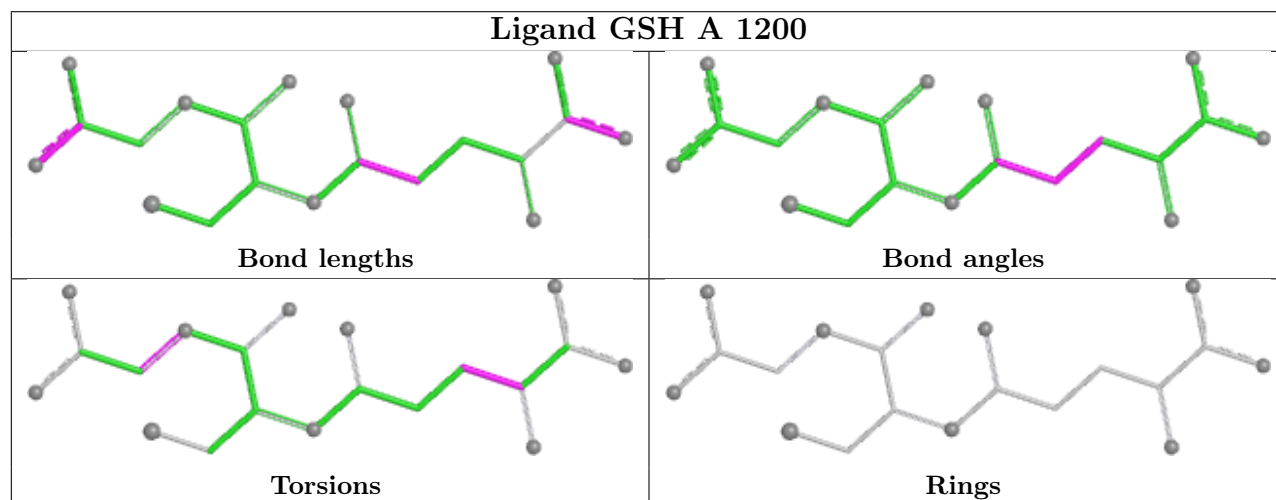
Mol	Chain	Res	Type	Atoms
3	A	1200	GSH	N1-CA1-CB1-CG1
3	C	1600	GSH	C1-CA1-CB1-CG1
3	D	1800	GSH	N1-CA1-CB1-CG1
3	D	1800	GSH	C1-CA1-CB1-CG1
3	C	1600	GSH	O2-C2-N3-CA3
3	B	1400	GSH	O2-C2-N3-CA3
3	D	1800	GSH	O2-C2-N3-CA3
3	D	1800	GSH	N2-CA2-CB2-SG2
3	A	1200	GSH	C1-CA1-CB1-CG1
3	B	1400	GSH	C1-CA1-CB1-CG1
3	B	1400	GSH	N1-CA1-CB1-CG1
3	D	1800	GSH	C2-CA2-CB2-SG2
3	A	1200	GSH	C3-CA3-N3-C2
3	C	1600	GSH	O2-C2-CA2-N2
3	C	1600	GSH	OE1-CD1-CG1-CB1
3	D	1800	GSH	N2-CD1-CG1-CB1
3	C	1600	GSH	N1-CA1-CB1-CG1

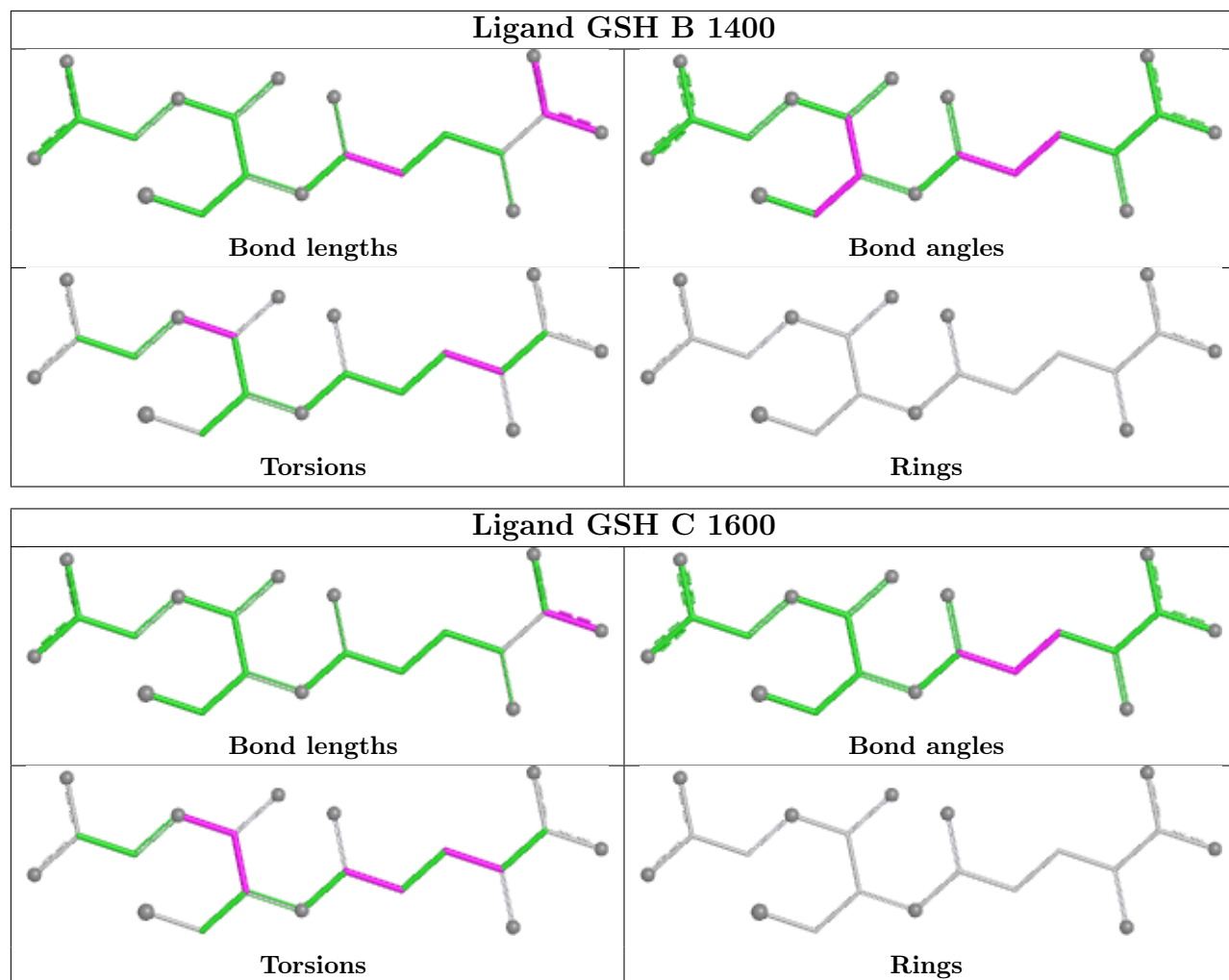
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1600	GSH	2	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	198/198 (100%)	0.34	6 (3%) 52 56	12, 20, 30, 37	0
1	B	198/198 (100%)	0.21	2 (1%) 79 82	11, 19, 27, 35	0
1	C	198/198 (100%)	0.53	10 (5%) 33 37	14, 22, 37, 40	0
1	D	198/198 (100%)	0.86	17 (8%) 16 17	14, 25, 41, 44	0
All	All	792/792 (100%)	0.49	35 (4%) 39 42	11, 21, 37, 44	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	722	TYR	4.1
1	A	110	ASP	4.1
1	D	710	ASP	3.6
1	A	144	MET	3.4
1	B	344	MET	3.4
1	D	634	ILE	3.4
1	D	799	LEU	3.3
1	A	145	SER	3.1
1	D	637	ALA	3.1
1	C	478	ALA	3.0
1	D	711	VAL	2.7
1	C	522	TYR	2.7
1	D	636	GLN	2.7
1	D	602	PRO	2.7
1	C	593	ARG	2.6
1	C	435	GLU	2.6
1	C	438	ASP	2.4
1	D	790	TRP	2.4
1	D	676	ASP	2.4
1	C	441	GLU	2.4
1	A	35	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	642	ILE	2.3
1	C	510	ASP	2.3
1	D	715	MET	2.3
1	C	517	ASN	2.3
1	D	793	ARG	2.2
1	D	639	TRP	2.2
1	B	310	ASP	2.2
1	D	623	ALA	2.1
1	C	544	MET	2.1
1	A	111	VAL	2.1
1	C	440	PRO	2.1
1	A	105	ALA	2.0
1	D	645	THR	2.0
1	D	674	ASN	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.

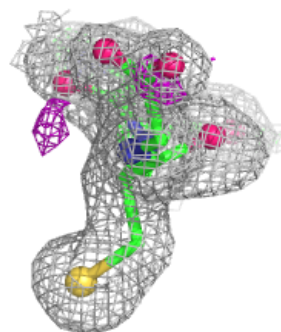
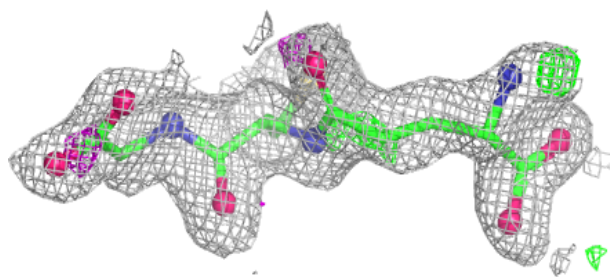
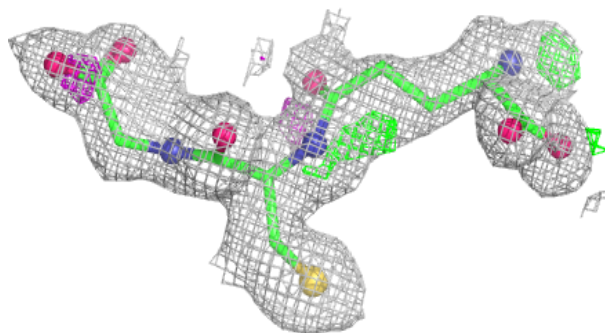
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GSH	D	1800	20/20	0.86	0.12	26,28,31,31	0
3	GSH	C	1600	20/20	0.88	0.11	26,27,29,29	0
3	GSH	A	1200	20/20	0.90	0.10	20,23,28,28	0
3	GSH	B	1400	20/20	0.93	0.09	18,20,23,23	0
2	MG	A	1901	1/1	0.98	0.04	22,22,22,22	0
2	MG	B	1902	1/1	0.99	0.07	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.

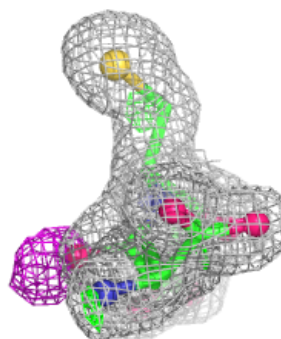
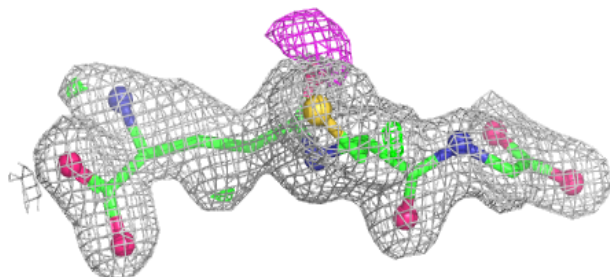
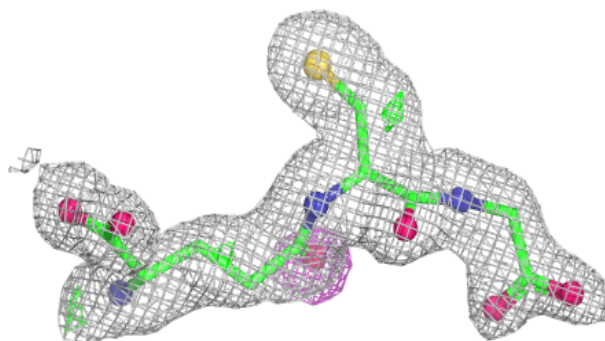
Electron density around GSH D 1800:

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



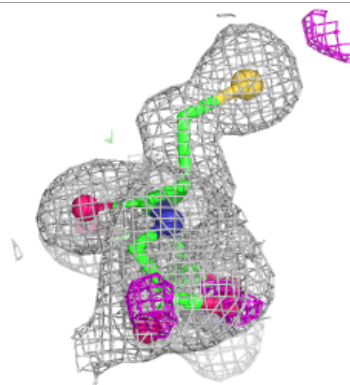
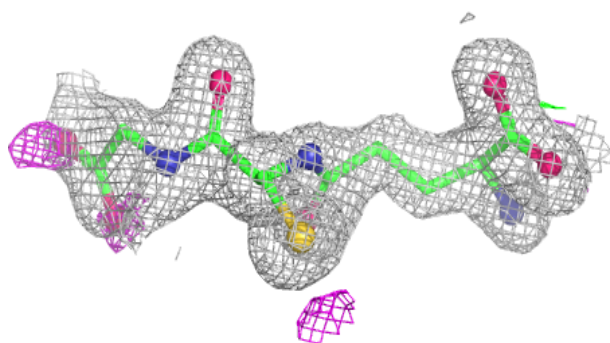
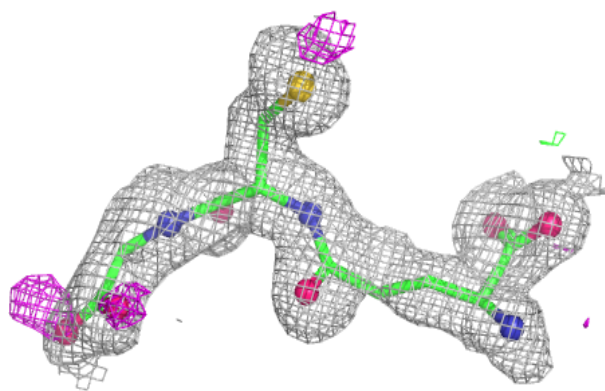
Electron density around GSH C 1600:

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

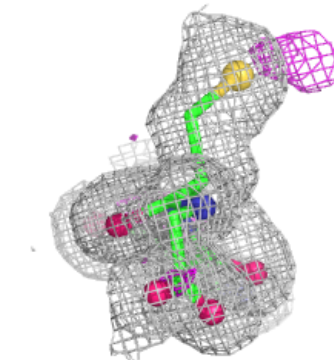
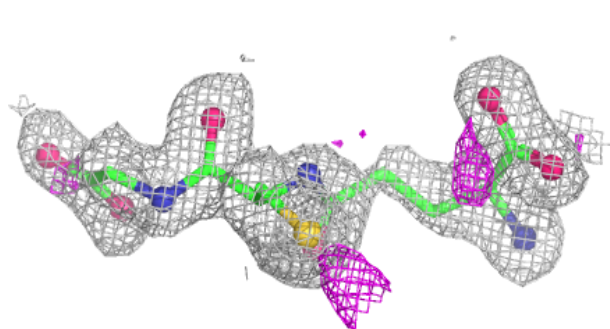
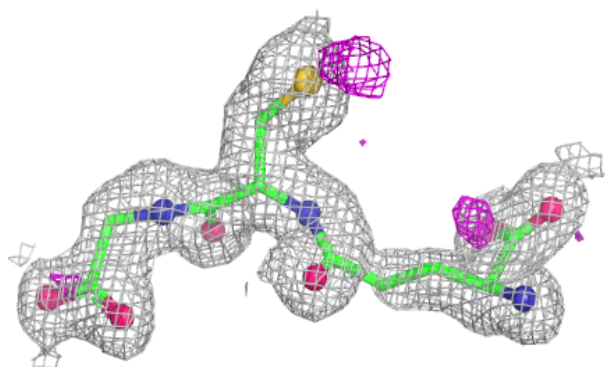


Electron density around GSH A 1200:

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

**Electron density around GSH B 1400:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.