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#### **Supplemental Information**

### **Structures of Human Antibodies Bound**

#### to SARS-CoV-2 Spike Reveal Common Epitopes

#### and Recurrent Features of Antibodies

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### Supplemental Tables

## Table S1. Crystallographic data collection and refinement statistics for C105 Fab

#### structure (related to Figure 5).

|   | C105 Fab                  |  |  |
|---|---------------------------|--|--|
|   | (12-1, SSRL)              |  |  |
| PDB ID  | 6XCA                      |  |  |
| Data collection <sup>a</sup>                            |                           |  |  |
| Space group   | 1222                      |  |  |
| Unit coll (Å)   | 67.4 120.1 122.2          |  |  |
|   | 07.4, 120.1, 123.5        |  |  |
| $(\alpha, \beta, \gamma(\beta))$                        | 1.0                       |  |  |
| $\mathbf{P}_{\text{assolution}}\left(\mathbf{A}\right)$ | 1.0                       |  |  |
| Lucique Deflections                                     | <i>46</i> 712 (2752)      |  |  |
| Completeness (%)  | 40,715(2752)<br>100(00.8) |  |  |
| Redundency  | 100(99.8)                 |  |  |
|   | 0.8 (0.3)                 |  |  |
| $CC_{1/2}$ (%)  | 98.8 (54.1)               |  |  |
| < <u>I</u> / <u></u> σI>                                | 5.7 (1.2)                 |  |  |
| Mosaicity (1)   | 0.19                      |  |  |
| R <sub>merge</sub> (%)                                  | 18.1 (157)                |  |  |
| R <sub>pim</sub> (%)                                    | 7.9 (70.5)                |  |  |
| Wilson B-factor   | 16.8                      |  |  |
|   |                           |  |  |
|   |                           |  |  |
| Refinement and Validation                               |                           |  |  |
| Resolution (Å)  | 38.9 - 1.80               |  |  |
| Number of atoms   |                           |  |  |
| Protein   | 3,132                     |  |  |
| Ligand  | 10                        |  |  |
| Waters  | 477                       |  |  |
| $R_{work}/R_{free}$ (%)                                 | 18.7/21.6                 |  |  |
| R.m.s. deviations                                       |                           |  |  |
| Bond lengths (Å)  | 0.006                     |  |  |
| Bond angles (°)   | 0.853                     |  |  |
| MolProbity score  | 1.29                      |  |  |
| Clashscore (all atom)                                   | 4.2                       |  |  |
| Poor rotamers (%)                                       | 0                         |  |  |
| Ramachandran plot                                       |                           |  |  |
| Favored (%)   | 97.6                      |  |  |
| Allowed (%)   | 2.4                       |  |  |
| Disallowed (%)  | 0                         |  |  |
| Average <i>B</i> -factor (Å)                            | 27.1                      |  |  |

<sup>a</sup>Numbers in parentheses correspond to the highest resolution shell

# Table S2. Cryo-EM data collection and refinement statistics for C105-S complex structure

(related to Figure 5).

|  | C105                                     | C105                                     |
|--|--|--|
|  | SARS-CoV-2 S 2P                          | SARS-CoV-2 S 2P                          |
|  | (state 1)                                | (state 2)                                |
| PDB  | 6XCM                                     | 6XCN                                     |
| EMD  | 22127                                    | 22128                                    |
| Microscope                                 | Titan Krios                              | Titan Krios                              |
| Camera                                     | Gatan K3 Summit                          | Gatan K3 Summit                          |
| Magnification                              | 105,000x                                 | 105,000x                                 |
| Voltage (kV)                               | 300                                      | 300                                      |
| Recording mode                             | counting                                 | counting                                 |
| Dose rate (e/pixel/s)                      | 22.1                                     | 22.1                                     |
| Electron dose $(e/Å^2)$                    | 60                                       | 60                                       |
| Defocus range (µm)                         | 1.0 - 2.5                                | 1.0 - 2.5                                |
| Pixel size (Å)                             | 0.418 (super resolution); 0.836 (binned) | 0.418 (super resolution); 0.836 (binned) |
| Micrographs collected                      | 5,940                                    | 5,940                                    |
| Micrographs used                           | 5,336                                    | 5,336                                    |
| Total extracted particles                  | 71,289                                   | 71,289                                   |
| Refined particles                          | 57,710                                   | 14,119                                   |
| Symmetry imposed<br>Nominal Resolution (Å) | C1                                       | C3                                       |
| FSC 0.5 (unmasked/masked)                  | 3.90/3.60                                | 4.00/3.60                                |
| FSC 0.143 (unmasked.masked)                | 3.40/3.20                                | 3.50/3.40                                |
| Map sharpening $B$ -factor                 |  |  |
| <b>Refinement and Validation</b>           |  |  |
| Number of atoms                            |  |  |
| Protein                                    | 25,973                                   | 27,798                                   |
| Ligand                                     | 711                                      | 873                                      |
| MapCC (global/local)                       | 0.86/0.84                                | 0.87/0.85                                |
| R.m.s. deviations                          |  |  |
| Bond lengths (Å)                           | 0.008                                    | 0.011                                    |
| Bond angles (°)                            | 0.812                                    | 0.893                                    |
| MolProbity score                           | 2.17                                     | 2.26                                     |
| Clashscore (all atom)                      | 13.6                                     | 14.8                                     |
| Poor rotamers (%)                          | 0.04                                     | 1.1                                      |
| Ramachandran plot                          |  |  |
| Favored (%)                                | 90.9                                     | 89.9                                     |
| Allowed (%)                                | 9  | 10.1                                     |
| Disallowed (%)                             | 0.1                                      | 0  |

Table S3. S protein mutations found in different SARS-CoV-2 isolates (related to Figure

**6)**.

| Mutation | Count | Frequency(%) | Location            |
|----------|-------|--------------|---------------------|
| D614G    | 9688  | 63.2         | S1 domain D         |
| P1263L   | 115   | 0.7          | S2 cytoplasmic tail |
| L5F      | 91    | 0.6          | signal sequence     |
| D936Y    | 88    | 0.6          | S2 HR1              |
| L54F     | 58    | 0.4          | S1 domain A         |
| G1124V   | 56    | 0.4          | S2                  |
| N439K    | 38    | 0.2          | S1 domain B (RBD)   |
| H49Y     | 35    | 0.2          | S1 domain A         |
| L18F     | 31    | 0.2          | S1 domain A         |
| L8V      | 30    | 0.2          | signal sequence     |
| A831V    | 29    | 0.2          | S2                  |
| D839Y    | 28    | 0.2          | S2                  |
| V483A    | 28    | 0.2          | S1 domain B (RBD)   |
| Q675H    | 24    | 0.2          | S1 domain D         |
| S50L     | 24    | 0.2          | S1 domain A         |
| S943P    | 22    | 0.1          | S2 HR1              |
| A1078S   | 21    | 0.1          | S2                  |
| R21I     | 19    | 0.1          | S1 domain A         |
| V367F    | 18    | 0.1          | S1 domain B (RBD)   |
| T29I     | 18    | 0.1          | S1 domain A         |

List of SARS-CoV-2 spike mutations with a frequency  $\geq 0.1\%$  in a set of 15335 isolates downloaded from the Global Initiative for Sharing All Influenza Data (GISAID) SARS-CoV-2 sequence database on 5/3/20 (Elbe and Buckland-Merrett, 2017; Shu and McCauley, 2017). The genomes were processed with the nextstrain augur pipeline (https://github.com/nextstrain/augur) (Hadfield et al., 2018), using MAFFT v7.464 (Katoh and Standley, 2013) for sequence alignment and FastTree (Price et al., 2010) to generate a phylogenetic tree. The resulting data were then analyzed with a custom Swift program.